

Dephasing and leakage dynamics of noisy Majorana-based qubits: Topological versus Andreev

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Topological quantum computation encodes quantum information nonlocally by nucleating non-Abelian anyons separated by distances L , typically spanning the qubit device size. This nonlocality renders topological qubits exponentially immune to dephasing from *all* sources of classical noise with operator support local on the scale of L . We perform detailed analytical and numerical analyses of a time-domain Ramsey-type protocol for noisy Majorana-based qubits that is designed to validate this coveted topological protection in near-term devices such as the so-called “tetron” design. By assessing dependence of dephasing times on tunable parameters, e.g., magnetic field, our proposed protocol can clearly distinguish a bona fide Majorana qubit from one constructed from semilocal Andreev bound states, which can otherwise closely mimic the true topological scenario in local probes. In addition, we analyze leakage of the qubit out of its low-energy manifold due to classical-noise-induced generation of quasiparticle excitations; leakage limits the qubit lifetime when the bulk gap collapses, and hence our protocol further reveals the onset of a topological phase transition. This experiment requires measurement of two nearby Majorana modes for both initialization and readout—achievable, for example, by tunnel coupling to a nearby quantum dot—but no further Majorana manipulations, and thus constitutes an enticing prebraiding experiment. Along the way, we address conceptual subtleties encountered when discussing dephasing and leakage in the context of Majorana qubits.

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I. INTRODUCTION

Developing a scalable quantum computing architecture that can withstand decoherence to the extent necessary for real-world applications poses an enormous scientific and technological undertaking. One path forward pursues Majorana-based topological qubits. These promise to provide robust protection against decoherence at the hardware level by storing quantum information nonlocally, specifically in a fermion-parity degree of freedom in a topological phase [1,2]. Since 2012, remarkable experimental progress [3–16] (see Ref. [17] for a recent review) has established the potential existence of Majorana zero modes in devices following the popular recipe [18,19] of engineering topological superconductivity [20] by interfacing s -wave superconductors with spin-orbit-coupled semiconducting nanowires subjected to modest magnetic fields.

Most of these laboratory efforts have, however, focused on end-of-wire tunneling conductance spectroscopy, a local measurement which may have difficulty differentiating bona fide well-separated Majorana zero modes in a topological phase [1] from near-zero-energy Andreev bound states (ABSs) arising in a trivial phase. One might naively expect that the latter should not “stick” near zero energy upon variation of system parameters such as magnetic field, chemical potential, etc. Important theoretical and numerical work [21–33] has nevertheless shown that seemingly quite reasonable spatial variations

in the potentials can trap ABSs—even deep in the trivial phase—that exhibit a high propensity to reside near zero energy (at least within the energy resolution of transport experiments). Andreev states, i.e., ordinary (complex) fermionic modes, can always be described in terms of two Majorana operators. Oftentimes, near-zero-energy ABSs are *semilocal* in that they consist of two Majorana operators separated in space by a finite distance, albeit one much less than the physical wire length [25,28,30,31,33]; the corresponding coupling strength between the constituent Majorana operators can thus be very small. This property endows semilocal ABSs with the (rather unfortunate) ability to masquerade as true Majorana zero modes even with respect to more detailed local-probe characteristics such as $2e^2/h$ zero-bias-peak conductance quantization [26,28] and 4π Josephson periodicity [34]. ABSs induced by local potential inhomogeneities arguably describe much of the experimental data as compellingly as true Majorana zero modes. While states encoded through semilocal ABSs might still furnish qubits with some degree of protection [30], such a situation is clearly suboptimal.

Various schemes have been proposed to distinguish Majorana zero modes from Andreev levels: (1) Detecting the bulk topological phase transition accompanying the onset of Majorana modes, e.g., through bulk-gap closure and reopening [35,36], surely provides an unambiguous fingerprint; see Ref. [37] for an experiment in this direction. Recently studied

multiterminal N-S-N devices provide an appealing platform for revealing the bulk phase transition via nonlocal conductance measurements [38–41]. (2) Majorana modes generate correlations in conductance measured at the two ends of a wire, whereas ABSs localized to each end generically produce no such correlations [25]. (3) Sensitivity of the energy to local perturbations, such as a quantum dot [13,42,43] or sharp edge potential [29], is expected to differ sharply for ABSs and Majorana modes due to the nonlocal, and hence energetically rigid, nature of the latter (see also Ref. [44]). (4) The current-phase relation in a Cooper-pair transistor has been shown to also differ qualitatively for these two scenarios [45]. (5) On the more challenging end, demonstration of inherently robust non-Abelian braiding operations is possible only with true Majorana zero modes.

We add to this list by analyzing in detail a time-domain protocol that not only unambiguously differentiates ABSs from Majorana modes, but further reveals important device characteristics (notably qubit lifetimes and precise Majorana-hybridization energies), exposes the topological phase transition, and clearly elucidates the enormous benefit of Majorana-based qubits insofar as robustly *protecting* quantum information. The main idea is to use *noise sensitivity* of a prototype qubit to validate (or invalidate) its topological nature. Similar ideas were originally put forth in Sec. IV of Ref. [46] in a related context. As we will see, in our framework ABSs that arise over a restricted parameter regime before the onset of well-separated Majorana modes in fact become a feature rather than a bug; their presence allows one to benchmark the quality of a bona fide topological qubit within a single device.

The remainder of this paper is organized as follows. Section II provides an executive summary of our main results. Section III introduces the microscopic models that we employ and also outlines our numerical simulation strategy. The Ramsey-type protocol that we use to probe the qubit dynamics is detailed in Sec. IV. Section V analytically estimates dephasing and leakage times for the qubit, while Sec. VI extracts these quantities from explicit time-dependent numerical simulations. We conclude with a discussion in Sec. VIII. Appendices provide supplemental calculational details.

II. EXECUTIVE SUMMARY

For concreteness, we focus on the so-called “tetron” Majorana qubit design shown in Fig. 1 [47–49] (equivalent for our purposes to the “loop qubit” [49]). This prototype topological qubit consists of two parallel semiconducting nanowires (gray) each proximitized by an *s*-wave superconductor and tuned to the topological phase with an external magnetic field (orange denotes topological superconductors). The concomitant four Majorana zero modes—defined through *instantaneous* eigenstates of the microscopic, time-dependent Hamiltonian—are denoted $\gamma_{1,2,3,4}$. A trivial superconductor (blue) bridges the two wires. In a real physical implementation, the entire mesoscopic device is floating so as to protect against quasiparticle poisoning events from the outside via charging energy [48]. For the nontopological ABS qubit¹

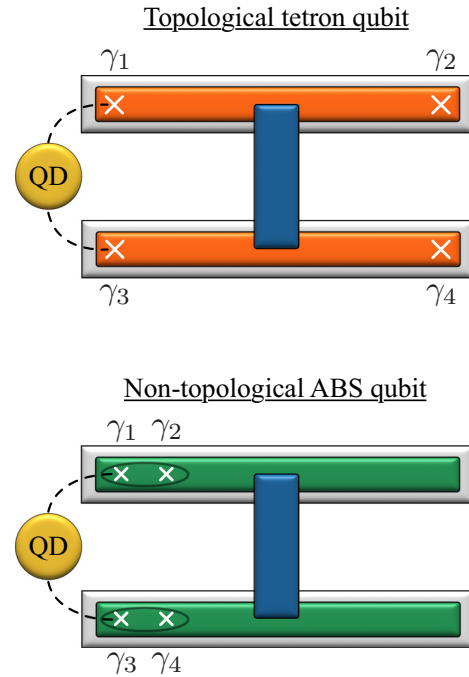


FIG. 1. Schematics of the topological tetron qubit (top) and a nontopological qubit arising from Andreev bound states (ABSs) residing near zero energy (bottom). Both cases consist of two parallel semiconducting nanowires (gray) proximity coupled to *s*-wave superconductors (orange and green, in the topological and nontopological scenarios, respectively); a trivial superconducting bridge connects the two parallel superconducting segments. As described in the text, the spatial separation of the low-energy Majorana modes $\gamma_{1,2,3,4}$ is qualitatively distinct in the two scenarios. The two leftmost such modes, i.e., γ_1 and γ_3 , are coupled to a nearby quantum dot (QD), allowing initialization and readout of the respective Majorana parity $i\gamma_1\gamma_3 = \pm 1$.

that we wish to contrast against, $\gamma_{1,2,3,4}$ instead represent a maximally localized set of Majorana operators that weakly hybridize at the left end of the device as illustrated in the bottom of Fig. 1. Unless specified otherwise, our discussion below pertains to both types of qubits.

Assuming that we are working on timescales much less than the characteristic poisoning time, and additionally that the system is confined to the low-energy, nearly degenerate ground-state manifold with high probability—e.g., the temperature and characteristic noise frequencies are well below the bulk excitation gap (see below)—we can encode logical qubit states by $|0\rangle \equiv |i\gamma_1\gamma_2 = +1\rangle$ and $|1\rangle \equiv |i\gamma_1\gamma_2 = -1\rangle$. [Here global fermion parity is fixed to $(i\gamma_1\gamma_2)(i\gamma_3\gamma_4) = +1$, so that specifying $i\gamma_1\gamma_2$ automatically specifies $i\gamma_3\gamma_4$.] In this basis, we can identify Pauli operators $Z \equiv i\gamma_1\gamma_2$, $X \equiv i\gamma_1\gamma_3$, and $Y \equiv -i\gamma_1\gamma_4$. Suppose that γ_1 and γ_2 couple with hybridization energy $\varepsilon_{12}(t)$ that is time dependent due to noise, and that γ_3 and γ_4 similarly couple with energy $\varepsilon_{34}(t)$. The time-averaged couplings decay asymptotically with the intrawire Majorana separation as an oscillatory exponential: $\varepsilon_{12} \sim \cos(k_{\text{top}}L_{\text{top}})e^{-L_{\text{top}}/\xi_{\text{top}}}$ and $\varepsilon_{34} \sim \cos(k_{\text{bot}}L_{\text{bot}})e^{-L_{\text{bot}}/\xi_{\text{bot}}}$, where k is related to the Fermi wave vector of the wire, ξ is the effective superconducting coherence length [50], and

¹For more parallel terminology, one can view the ABS qubit as a superficial, hardly topological tetron (SHT tetron).

“top”/“bot” refers to the top and bottom wires of the tetron. For the topological qubit, $L_{\text{top/bot}}$ are given by the wire length L —thus taking full advantage of the exponential suppression. By contrast, for the ABS qubit $L_{\text{top/bot}}$ is on the scale of the coherence length or smaller, and hence the splitting need not conform to such an exponential. In principle, couplings between Majoranas on different wires, such as ε_{14} , ε_{23} , etc., can also be present. However, since the separation between these Majorana modes generally exceeds that between Majorana modes on the same wire, these couplings will be significantly smaller than the intrawire hybridization energies—at least if the superconducting bridge connecting the two wires is sufficiently long and well gapped. In this case, their primary effect is to reduce the amplitude of qubit oscillations and slightly shift the qubit precession frequency while leaving the qualitative features unaffected. We can therefore safely ignore such subdominant terms and write the Hamiltonian governing the dynamics of the qubit as

$$\begin{aligned} H(t) &= \frac{i}{2} [\varepsilon_{12}(t) \gamma_1 \gamma_2 + \varepsilon_{34}(t) \gamma_3 \gamma_4] \\ &= \frac{1}{2} [\varepsilon_{12}(t) + \varepsilon_{34}(t)] Z \equiv \frac{1}{2} E(t) Z. \end{aligned} \quad (1)$$

Finally, a quantum dot (ignored thus far) sits proximate to the left ends of the nanowires, thereby allowing a joint parity readout [48,51,52] of nearby Majoranas γ_1 and γ_3 , i.e., the Pauli X operator.

Our proposed protocol is conceptually very simple: (1) Initialize the system into an X eigenstate by measuring $i\gamma_1\gamma_3$; (2) let the system evolve unitarily for a wait time t ; and finally (3) remeasure X . Assuming the unitary evolution is governed by Eq. (1), which ignores interwire Majorana hybridization as well as higher-energy excitations, the qubit precesses about the z axis on the equator of the Bloch sphere. Let us write the instantaneous qubit splitting defined through Eq. (1) as $E(t) = \hbar\omega_0 + \delta E(t)$. Here, ω_0 is the *time-averaged* qubit precession frequency while $\delta E(t)$ encodes the effects of classical noise and is responsible for dephasing. Note that we assume here that $\delta E(t)$ fluctuates around zero mean. Taking the initial state to be an eigenstate of $X = i\gamma_1\gamma_3$ with eigenvalue $+1$, noise averaging the X readout measurement gives

$$\langle \psi(t) | X | \psi(t) \rangle = \cos(\omega_0 t) f(t/T_2), \quad (2)$$

where the envelope function $f(t/T_2)$ decays on a timescale given by the *dephasing time* T_2 . This experiment therefore simultaneously probes ω_0 (and thus reveals the mean Majorana hybridization energy) as well as T_2 —both critical device characteristics that differ starkly for topological and ABS qubits. We note that the described protocol is similar in spirit to the so-called Ramsey sequence [53,54]. The main differences are that we (1) do not include an external driving field transverse to Z and (2) initialize and read out the qubit with projective X measurements in contrast to utilizing $\pi/2$ pulses and Z initialization/readout²

²A related sequence involving Z initialization, two $\pi/2$ rotations about the x axis buttressing a wait time t , followed by a final Z measurement was proposed in Ref. [46], as appropriate for the qubit design presented therein.

A particularly advantageous feature of Majorana-based topological qubits is that, due to their inherently nonlocal encoding, the splitting E remains exponentially small—and thus also “exponentially flat”—in response to changes in *all* local Hamiltonian couplings (e.g., chemical potential, Zeeman field, etc.). This property endows topological qubits with exponential protection from all classical noise sources, yielding the following nontrivial scaling relation that connects time-averaged splitting and dephasing [46]:

$$\frac{1}{\omega_0} \sim T_2^a \sim e^{L/\xi}. \quad (3)$$

Here, a is an order-one number dependent on details of the noise. In other words, as the topological qubit becomes “perfect,” the splitting vanishes and the dephasing time diverges in a correlated fashion. For comparison, the transmon qubit [55] benefits similarly from an exponential protection from charge noise, but not other noise sources; additionally, there the qubit splitting need not be exponentially small. The prebraiding protocol sketched above and analyzed in detail below is capable of probing the scaling relation in Eq. (3), which can for practical purposes be taken as a *definition* of a topologically protected quantum memory. Note also that as the qubit quality improves (i.e., as L/ξ increases) the splitting $\hbar\omega_0$ becomes more difficult to resolve in transport [9,56] yet easier to resolve in the time-domain measurements employed in our protocol.

The trivial ABS qubit exhibits qualitatively different behavior. First, since the requisite “accidental” low-energy Andreev bound states would emerge due to some local, nonuniversal features in the potential landscapes near the ends of the wires, the ABS qubit would clearly not exhibit exponential protection against arbitrary local noise sources. Moreover, neither ω_0 nor T_2 will vary appreciably with wire length L , in sharp contrast to the scaling relation in Eq. (3) that uniquely identifies the topological qubit.

While it would be ideal to perform experiments at different L , keeping all other parameters fixed in this process has its practical challenges. We nevertheless argue that one can compellingly distinguish the two scenarios in a *single* prototype qubit at fixed L by carrying out the protocol at different field strengths B . Numerous numerical simulations [24,25,28,30,31,33] indicate a tendency for low-energy Andreev bound states to form over an extended field interval below the onset of a topological phase hosting true Majorana zero modes. In such a scenario, upon increasing B from zero, the system first realizes an ABS qubit, then encounters a topological phase transition at a critical field B_c , and finally forms a topological qubit.

The behavior of ω_0 and T_2 during this evolution reveals a wealth of information. In the ABS regime there is likely no clear universal behavior present, though the measured coherence times would provide a useful baseline. When B approaches B_c , the bulk gap becomes close to zero. Here, even low-frequency noise can efficiently excite the qubit out of the computational subspace, causing a precipitous reduction in the qubit lifetime. (Below we denote the characteristic “leakage” time associated with such excitations by T_{leak} .) Hence, our protocol provides a means of detecting the topological phase transition. As B increases beyond B_c , the gap reopening

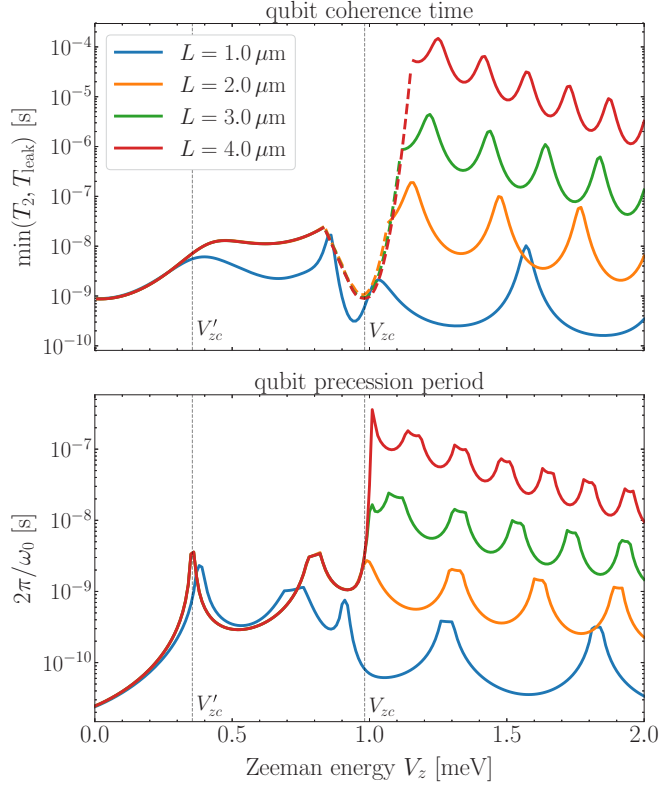


FIG. 2. Qubit coherence time defined as $\min(T_2, T_{\text{leak}})$ (top) and qubit precession period $2\pi/\omega_0$ (bottom) as a function of Zeeman energy V_z for tetron qubits constructed from the spinful nanowire model of Eq. (5) at different wire lengths L . For $V'_z < V_z < V_{zc}$, the qubit is of the nontopological ABS variety (Fig. 1, bottom panel), while for $V_z > V_{zc}$, the qubit is topological (Fig. 1, top panel). In the top panel, solid curves indicate that the qubit lifetime is *dephasing limited* ($T_2 < T_{\text{leak}}$), while dashed curves indicate that the qubit lifetime is *leakage limited* ($T_{\text{leak}} < T_2$). These calculations were performed using the quasianalytic methods of Sec. V; for more details, including the specific parameters chosen, please see Sec. VIB.

and concomitant appearance of robust Majorana zero modes rapidly boost the qubit's coherence time—ideally to values exceeding the ABS-qubit lifetime by orders of magnitude. The scaling relation in Eq. (3), which should now be viewed in terms of field-induced variation of ξ at fixed L , becomes operative in this topological regime. The topological-qubit frequency ω_0 additionally exhibits characteristic oscillations with magnetic field, reflecting oscillatory overlap of the Majorana wave functions [50], in turn yielding out-of-phase oscillations in the dephasing time T_2 [46]. Further increasing B eventually suppresses the gap for the topological phase and hence increases ξ , thereby diminishing $1/\omega_0$ and T_2 in accordance with Eq. (3). Note also that for sufficiently long topological wires, the exponentially long dephasing time T_2 is eventually cut off by the leakage time T_{leak} (which is set by the L -independent excitation gap).

The above points are demonstrated in Fig. 2, where we show concrete calculations of the qubit coherence time defined as $\min(T_2, T_{\text{leak}})$ and qubit precession period $2\pi/\omega_0$ for

noisy tetron qubits modeled within the canonical single-band description of proximitized spin-orbit-coupled nanowires [18,19]. Here, the horizontal axis denotes the Zeeman energy, related to the magnetic field B through $V_z = \frac{1}{2}g\mu_B B$, with g the effective Landé g factor (assumed equivalent for both wires of the tetron) and μ_B the Bohr magneton. In these plots, and in all simulations in this paper, we neglect dependence of the noise on magnetic field and other parameters. We stress that although this assumption will undoubtedly be violated to some degree in experiment, the nonmonotonicity and sheer magnitude of the effects illustrated in Fig. 2 are unlikely to be modified significantly by such nonuniversal noise properties. For more details of the model Hamiltonian, means of calculation, and chosen physical parameters for the data in Fig. 2, see Secs. III A, V, and VIB below.

Many previous studies have explored the influence of noise on Majorana-based qubits from various perspectives (see, e.g., Refs. [49,57–67]). We emphasize that here we are proposing to use noise to our *advantage* in verifying the topological nature of a prototype Majorana qubit, and as a by-product are able to extract crucial device characteristics that would otherwise be difficult to obtain. Time-domain experiments of the type that we propose certainly entail significant challenges, but they are arguably a prerequisite for successful demonstration of exponentially accurate braiding gates.

III. MICROSCOPIC MODELS AND THEORETICAL FORMULATION

A. Microscopic Hamiltonians

We exploit two different microscopic Hamiltonians for the one-dimensional (1D) topological superconductors comprising our prototype tetron qubits. For a truly minimal description we appeal to Kitaev's model [1] for a 1D spinless p -wave superconductor, given by the Hamiltonian

$$H_K = \sum_{i=1}^L \left[-\mu c_i^\dagger c_i - \frac{1}{2} (J c_i^\dagger c_{i+1} + \Delta c_i c_{i+1} + \text{H.c.}) \right]. \quad (4)$$

Here, μ is the chemical potential, $J \geq 0$ is the nearest-neighbor hopping strength, and Δ is the p -wave pairing strength (assumed real). The topological phase occurs for $|\mu| < J$, wherein a chain with open boundary conditions harbors a pair of Majorana zero modes (MZMs), one at each end.

For a more realistic treatment, we consider the one-band spinful model of a Rashba spin-orbit-coupled nanowire proximitized with an s -wave superconductor and subjected to a Zeeman field [18,19]. The Hamiltonian reads as

$$H_{\text{NW}} = \int_0^L dx \left[\psi^\dagger \left(-\frac{\hbar^2 \partial_x^2}{2m} - \mu + V(x) - i\alpha \sigma^y \partial_x + V_z \sigma^z \right) \psi + \Delta(x) (\psi_\uparrow \psi_\downarrow + \text{H.c.}) \right], \quad (5)$$

where m is the effective electron mass, μ is the chemical potential, α is the Rashba spin-orbit strength, $V_z = \frac{1}{2}g\mu_B B$ is the Zeeman energy, $\Delta(x)$ is the induced s -wave pairing

amplitude, and Pauli matrices $\sigma^{x,z}$ act in spin space. Note that we explicitly included a spatially varying external electric potential $V(x)$ and also allowed the pairing potential $\Delta(x)$ to depend on position: these features are essential for creating conditions that trap near-zero-energy ABS's that mimic MZM's. A uniform system with $V(x) = 0$ and $\Delta(x) = \Delta_0$ resides in the topological phase provided $h > \sqrt{\Delta_0^2 + \mu^2}$. For numerical evaluation we discretize Eq. (5) in a standard way; when essential, we will specify the lattice spacing used in this discretization procedure below.

At fixed global fermion parity, we microscopically model tetron qubits via

$$H = H_{\text{top}} + H_{\text{bot}}, \quad (6)$$

where H_{top} and H_{bot} take the form of either Eqs. (4) or (5) and, respectively, describe the top and bottom topological superconductors in the qubit. Classical noise is readily included by endowing parameters in H with stochastic time dependence. In principle, one should include a “bridge” term that couples the two superconductors (see Fig. 1). We assume for simplicity that the bridge yields only small quantitative effects and thus ignore its presence in our calculations. Both of our microscopic models describe noninteracting electrons in a non-number-conserving formulation and thus fail to capture charging-energy effects that would be present in a laboratory realization of the tetron (see Sec. II). This simplification is assumed for obvious computational tractability reasons; however, the effects of $1/f$ charge noise, as was argued to be the dominant physical source of tetron dephasing in Ref. [49], can be captured at a rough qualitative level by taking the global chemical potentials to fluctuate stochastically in Eqs. (4) and (5).

B. Majorana-operator reformulation

In our microscopic numerical simulations of tetrans, we find it most convenient to work in the language of *local* Majorana operators a_i —with $\{a_i, a_j\} = 2\delta_{ij}$, $a_i^\dagger = a_i$, and $(a_i)^2 = 1$ —by decomposing the local Dirac fermions in the usual way, e.g., $c_i = (a_{2i-1} - ia_{2i})/2$ in the Kitaev model context [1]. In this basis, the (quadratic) microscopic tetron Hamiltonian is represented by an $N \times N$ antisymmetric matrix $A = -A^T$:

$$H = \frac{i}{4} \sum_{i,j=1}^N A_{ij} a_i a_j = \frac{i}{4} \mathbf{a}^T A \mathbf{a}, \quad (7)$$

with N the total number of local Majorana operators; for spinless (spin- $\frac{1}{2}$) models with N_{sites} physical sites, $N = 2N_{\text{sites}}$ ($N = 4N_{\text{sites}}$). It is a relatively straightforward numerical exercise (see, e.g., Refs. [68–70]) to bring A into so-called “canonical form” via an orthogonal rotation U :

$$B = U^T A U = \bigoplus_{k=1}^{N/2} \varepsilon_k i \sigma^y, \quad (8)$$

such that

$$H = \frac{i}{4} \mathbf{b}^T B \mathbf{b} = \frac{i}{2} \sum_{k=1}^{N/2} \varepsilon_k b_{2k-1} b_{2k}. \quad (9)$$

The canonical Majorana modes $b_i = [\mathbf{b}]_i$ are related to the original local Majoranas $a_i = [\mathbf{a}]_i$ via the orthogonal transformation $\mathbf{b} = U^T \mathbf{a}$, and $\varepsilon_k \geq 0$ are the energies (or instantaneous energies when the Hamiltonian depends on time).

Due to Wick's theorem, any quadratic system can be completely described by the real, antisymmetric covariance matrix

$$M_{ij} = \frac{-i}{2} \langle [a_i, a_j] \rangle, \quad (10)$$

where the square brackets denote a commutator and $\langle \dots \rangle = \text{Tr}\{\rho \dots\}$ a quantum expectation value taken with respect to density matrix ρ ; if $\rho = |\psi\rangle\langle\psi|$ represents a pure state, one finds $M^2 = -1$. As an example, the ground state of H corresponds to a covariance matrix

$$M_{\text{g.s.}} = U M_y U^T, \quad (11)$$

with

$$M_y = \bigotimes_{k=1}^{N/2} i \sigma^y \quad (12)$$

a reference covariance matrix in the canonical basis of the modes b_i which encodes $i b_{2k-1} b_{2k} = -1 \forall k = 1, \dots, N/2$. It is straightforward to construct more general states by appropriately toggling elements of M_y (see Refs. [69,70] and Sec. IV).

Finally, the Schrödinger equation for Gaussian states takes the form of the following ordinary matrix differential equation for the covariance matrix:

$$\frac{dM(t)}{dt} = [A(t), M(t)], \quad (13)$$

where $A(t)$ specifies the (possibly time-dependent) Hamiltonian [cf. Eq. (7)] and the square brackets denote the matrix commutator. Any physical observable can then be computed with knowledge of $M(t)$; relevant examples of such measurements will be described in the next section.

IV. RAMSEY-TYPE PROTOCOL

A. Qubit definition

In what follows, we assume for concreteness that the devices in Fig. 1 possess even global fermion parity. We will encode the qubit in the subspace consisting of the two lowest-lying *instantaneous* even-parity eigenstates of the microscopic Hamiltonian $H(t)$ that depends on time due to noise. These presumed nearly degenerate states could arise because the system either realizes bona fide topological phases yielding four well-separated Majorana modes (upper panel of Fig. 1), or exhibits a pair of “accidentally” low-energy Andreev bound states (lower panel of Fig. 1). Since we will explore platforms that contain both types of near degeneracies depending on parameters, it is useful to treat them in a common framework. To this end we introduce a quartet of *maximally localized* (in real-space) Majorana operators $\gamma_{1,2,3,4}(t)$ that span the instantaneous qubit subspace at time t . In the topological case, residual overlap between Majorana operators splits the degeneracy by an energy that is exponentially small in their separation; in the low-energy Andreev-bound-state scenario,

pairs of Majorana operators sit in close proximity yet happen to hybridize weakly.

For a more precise definition of the qubit, suppose that

$$P(t) \equiv [i\gamma_1(t)\gamma_2(t)][i\gamma_3(t)\gamma_4(t)] = +1 \quad (14)$$

in the lowest-lying even-fermion-parity states, and define Pauli operators

$$Z_t = i\gamma_1(t)\gamma_2(t), \quad X_t = i\gamma_1(t)\gamma_3(t). \quad (15)$$

Our logical qubit states are then the minimum-energy many-body states with $Z_t = \pm 1$, i.e.,

$$|0_t\rangle \equiv |Z_t = +1\rangle, \quad |1_t\rangle \equiv |Z_t = -1\rangle. \quad (16)$$

The subscript t (suppressed in Sec. II) is included here and below as a reminder that these states are defined with respect to Majorana operators extracted from the instantaneous Hamiltonian at time t , rather than with respect to a fixed basis.

Some discussion is warranted regarding the definition of the maximally localized Majorana operators $\gamma_{1,2,3,4}(t)$ in our microscopic numerical simulations. The canonical modes b_i of the quadratic Hamiltonian H defined above [cf. Eqs. (8) and (9)] are not unique since arbitrary $\text{SO}(2)$ rotations $O_k^{2 \times 2} = \exp(i\theta_k \sigma^y)$ within each 2×2 block of Eq. (8) leave the matrix B invariant:

$$B = O B O^T, \quad O = \bigoplus_{k=1}^{N/2} O_k^{2 \times 2}. \quad (17)$$

Thus, Majorana modes $\tilde{b}_i = [\tilde{\mathbf{b}}]_i$ defined through

$$\tilde{\mathbf{b}} = O^T \mathbf{b} = \tilde{U}^T \mathbf{a}, \quad (18)$$

with $\tilde{U} = U O$, also form a canonical set with the same instantaneous energies, i.e.,

$$H = \frac{i}{4} \tilde{\mathbf{b}}^T B \tilde{\mathbf{b}} = \frac{i}{2} \sum_{k=1}^{N/2} \varepsilon_k \tilde{b}_{2k-1} \tilde{b}_{2k}. \quad (19)$$

The Majorana modes $\gamma_{1,2,3,4}(t)$ depicted in Fig. 1 refer to $\tilde{b}_{1,2,3,4}$ with $O_{k=1,2}^{2 \times 2}$ chosen such that the associated wave functions maximally localize at a given time t . Specifically, we first find a set of canonical near-zero-energy modes $b_{1,2,3,4}$ by bringing A into canonical form using the software package presented in Ref. [68]; in the local basis of the a_i , these modes are represented by column vectors of the matrix U , say $U_{i,j=1,2,3,4}$. We then optimize the $O_{k=1,2}^{2 \times 2}$ (each characterized by an angle $0 \leq \theta_k < 2\pi$) such that for the resulting $\tilde{U} = U O$ the fourth moments $\sum_{j=1}^2 \sum_{i=1}^N |\tilde{U}_{ij}|^4$ and $\sum_{j=3}^4 \sum_{i=1}^N |\tilde{U}_{ij}|^4$ (for $k = 1, 2$, respectively) are maximized; now, in the local a_i basis, these maximally localized modes $\tilde{b}_{1,2,3,4}$ are given by column vectors of \tilde{U} , i.e., $\tilde{U}_{i,j=1,2,3,4}$. We order these modes according to their real-space locations as shown in Fig. 1, thereby giving the desired $\gamma_{1,2,3,4}(t)$.³

As an aside, the numerical procedure spelled out above is tailored to the situation where the two superconductors

comprising the qubit decouple, which again we assume for simplicity is the case throughout. The Majorana modes $\gamma_{1,2}(t)$ [$\gamma_{3,4}(t)$] then have support entirely on the top (bottom) wire of the tetron. In the more realistic scenario in which a nonzero “bridge” Hamiltonian couples the wires, a physically plausible definition of $\gamma_{1,2,3,4}(t)$ involves instead maximizing the single fourth moment $\sum_{j=1}^4 \sum_{i=1}^N |\tilde{U}_{ij}|^4$ via a single orthogonal rotation $O^{4 \times 4}$ acting on the $b_{1,2,3,4}$. In the latter case, $\gamma_{1,2,3,4}(t)$ are not generally eigenmodes of the Hamiltonian.

B. Protocol details

Section II summarized our time-domain Ramsey-type protocol for the physical implementation of the tetron. In the “ideal” case the protocol involves (1) initializing the system into, say, the $X_{t=0} = +1$ eigenstate at time $t = 0$ via energy-level spectroscopy on the nearby quantum dot which tunnel couples to γ_1 and γ_3 , (2) letting the system evolve freely for time t under the influence of classical noise, and (3) reading out $X_t = \pm 1$ with the same quantum dot used for initialization. (For practical purposes, we will sometimes depart from this ideal protocol by initializing and measuring with respect to a fixed basis.) In the remainder of this section, we elaborate on these steps to set the stage for our analytic treatment of dephasing and noise in Sec. V and our full time-dependent microscopic simulations of tetron qubits in Secs. VI and VII.

1. Initialization

We start with initialization via the proximate quantum dot. At the level of our analysis, we assume that the quantum dot is “perfect” in the sense that a measurement projects exactly and instantaneously onto the qubit state

$$|\psi(0)\rangle = \frac{1}{\sqrt{2}}(|0_{t=0}\rangle + |1_{t=0}\rangle) \quad (20)$$

with $X_{t=0} = +1$ at time $t = 0$. Under this highly idealized assumption, we need not explicitly include the quantum dot in our analysis. In our microscopic simulations of noisy tetron qubits in Sec. VI, we use the *time-averaged* Hamiltonian—specified by an antisymmetric matrix A_0 —to define the relevant $\gamma_{1,2,3,4}$ used for initialization, which we denote $\gamma_{1,2,3,4}^{(0)}$. [Obtaining the modes instead using the initial Hamiltonian $A(t=0) \neq A_0$ for each noise realization, as implied by Eq. (20), leads to only negligible quantitative differences in the results presented there; see Sec. VII for more discussion on this point.] Specifically, we bring A_0 into canonical form as described above in Sec. III [see Eq. (9)]; we subsequently find the corresponding set of maximally localized near-zero-energy Majorana modes, which are encoded in column vectors of an orthogonal matrix \tilde{U}_0 . Next, we construct a reference covariance matrix $M_y(t=0)$ [cf. Eqs. (11) and (12)] in the basis of these modes $\tilde{b}_i^{(0)}$ corresponding to, say, $X_0 = i\gamma_1^{(0)}\gamma_3^{(0)} = +1$ and $i\gamma_2^{(0)}\gamma_4^{(0)} = -1$. The global fermion parity thus reads as $(i\gamma_1^{(0)}\gamma_2^{(0)})(i\gamma_3^{(0)}\gamma_4^{(0)}) = +1$, which for all presented simulations coincides with the parity of the absolute ground state of A_0 . (Recall that modes $\tilde{b}_{i=1,2,3,4}$ within the low-energy manifold correspond to the maximally localized modes $\gamma_{1,2,3,4}$, while the remainder $\tilde{b}_{i \neq 1,2,3,4}$ are equivalent to the respective original canonical modes b_i .) Finally, the

³This method for obtaining maximally localized near-zero-energy Majorana modes working entirely in the local Majorana representation parallels that described in Ref. [25] using the more traditional Bogoliubov–de Gennes (BdG) framework.

initial covariance matrix in the a_i basis [cf. Eq. (10)] reads as $M(t=0) = \tilde{U}_0 M_y(t=0) \tilde{U}_0^T$.

2. Unitary evolution

Once initialized, the system undergoes unitary time evolution with respect to the noisy Hamiltonian $H(t)$. After time t the initial state becomes

$$\begin{aligned} |\psi(t)\rangle &= U(t) |\psi(0)\rangle \\ &\equiv a_t |0_t\rangle + b_t |1_t\rangle + c_t |v_t\rangle, \end{aligned} \quad (21)$$

where $U(t)$ is the time-evolution operator (we discuss the second line shortly). Our microscopic numerical simulations are instead carried out in the Heisenberg picture, for which unitary evolution is encoded in Eq. (13); there, noise is included through time dependence in $A(t)$, which is essentially the Hamiltonian expressed in the Majorana basis. Appendix A details our treatment of noise (e.g., the procedure we employ for generating individual noise trajectories for a given realization) and numerical solution of Eq. (13).

We stress that in general $|0_t\rangle \neq U(t) |0_{t=0}\rangle$ since $|0_t\rangle$ is defined through instantaneous $H(t)$ eigenstates (and similarly for $|1_t\rangle$).⁴ In the second line of Eq. (21) we thus expressed the time-evolved state in terms of qubit states $|0_t\rangle$ and $|1_t\rangle$ at time t , and a ket $|v_t\rangle$ that signifies dynamically generated excited states. Generation of weight on the latter via nonzero c_t corresponds to “leakage” of the qubit away from the computational subspace. Leakage can arise from an odd or even number of fermionic excitations, thus respectively flipping or preserving $P(t)$. Using Eq. (21) we can define a qubit leakage time T_{leak} by writing

$$\overline{|a_t|^2} + \overline{|b_t|^2} \equiv 1 - f(t/T_{\text{leak}}), \quad (22)$$

where the overline indicates noise averaging. On the right side, $f(t/T_{\text{leak}})$ is a system-dependent function that vanishes at $t=0$ and grows—thus shifting weight onto excited states—on a characteristic timescale T_{leak} . An alternative leakage metric can be extracted from

$$\overline{\langle \psi(t) | P(t) | \psi(t) \rangle} \equiv 1 - \tilde{f}(t/\tilde{T}_{\text{leak}}) \quad (23)$$

with \tilde{f} a different function that grows on a timescale \tilde{T}_{leak} . The latter measures the weight on excited states containing only odd numbers of fermionic excitations, and thus provides an upper bound on the leakage time defined above, i.e., $T_{\text{leak}} \leq \tilde{T}_{\text{leak}}$.

At this point it is worth commenting on an alternative scheme wherein one defines the qubit in terms of a fixed basis of maximally localized Majorana operators obtained from, say, the time-averaged Hamiltonian A_0 or, for a given noise realization, the initial Hamiltonian $A(t=0)$. For an extreme case, suppose that $H(t)$ supports exact instantaneous zero-energy Majorana modes at any time t , but that *adiabatic* noise causes the locations of the Majorana modes to vary with time. The system will then evolve out of the low-energy subspace

spanned by the fixed Majorana operators, whereas the qubit subspace should clearly be perfectly preserved—as captured by tracking instantaneous Majorana operators. In general we expect that employing a fixed basis *overestimates* qubit errors, with the difference being most pronounced when the noise is “slow” and of “large” amplitude. We will quantify such effects later in Sec. VII.

3. Readout

The final step of the protocol involves readout. We again assume that the nearby quantum dot can perform this task, i.e., measure $X_t = \pm 1$, instantaneously and without error. Under this assumption, the relevant physical quantity is the noise-averaged expectation value of X_t with respect to the time-evolved state:

$$\overline{\langle \psi(t) | X_t | \psi(t) \rangle} = \cos(\omega_0 t) f(t/T_2), \quad (24)$$

which defines the qubit precession frequency ω_0 and dephasing time T_2 . Compared to Eq. (2), the t subscript on the left side explicitly indicates that the expectation value is (ideally) taken with respect to the instantaneous Majoranas at time t .

To calculate the quantum expectation value $\langle \psi(t) | X | \psi(t) \rangle$ for a given noise realization in our microscopic simulations, we basically invert the initialization procedure described above. Specifically, we conjugate $M(t)$ with \tilde{U} (which defines the measurement basis; see below) to arrive at

$$M_y(t) = \tilde{U}^T M(t) \tilde{U}, \quad (25)$$

from which any $\langle i\gamma_i \gamma_j \rangle$ can be read off directly. Measurements involving a product of more than two operators can be evaluated with use of Wick’s theorem; an important example is $P(t)$, which (partially) probes leakage out of the computational subspace [recall Eq. (23)].

As with initialization, there exists a choice as to precisely what is meant by the Majorana operators $\gamma_{1,2,3,4}$ being measured. Namely, these could be defined as (maximally localized) near-zero-energy modes with respect to, say, the *fixed* time-averaged Hamiltonian A_0 or the *instantaneous* Hamiltonian $A(t)$.⁵ Operationally, to measure in the basis corresponding to the time-averaged Hamiltonian such that $\gamma_{1,2,3,4} = \gamma_{1,2,3,4}^{(0)}$, we use Eq. (25) with $\tilde{U} = \tilde{U}_0$ (the same orthogonal matrix used in the initialization procedure). On the other hand, to measure in the instantaneous basis we need to find a set of maximally localized modes $\gamma_{1,2,3,4} = \gamma_{1,2,3,4}(t)$ at every time t , which are encoded in a time-dependent orthogonal matrix $\tilde{U} = \tilde{U}(t)$ to be used in Eq. (25). In our microscopic simulations, we implement both instantaneous and fixed measurement bases. Employing a fixed basis is computationally cheaper as it does not require bringing the Hamiltonian into canonical form and finding maximally localized modes at each time step. Additionally, there is an inherent sign ambiguity in the canonical modes (and thus also maximally localized modes) upon bringing $A(t)$ into canonical form numerically, i.e., the Hamiltonian is invariant

⁴ $|0_t\rangle$ and $|1_t\rangle$ are the two lowest-energy eigenstates of the instantaneous Hamiltonian in the same global parity sector as the evolving state $|\psi(t)\rangle$ (assumed even in this discussion), which may or may not coincide with the parity of the absolute instantaneous ground state.

⁵An alternative choice similar to the former case could be to use modes derived from the fixed *initial* Hamiltonian $A(t=0) \neq A_0$ for a given noise realization. We take this approach in Sec. VII.

upon taking $b_{2k-1} \rightarrow -b_{2k-1}$ and $b_{2k} \rightarrow -b_{2k}$ for a given k in Eq. (9). Therefore, measurements in the instantaneous basis such as $\langle i\gamma_1(t)\gamma_3(t) \rangle$ will be plagued by $\gamma_1(t)$ and $\gamma_3(t)$ having arbitrary relative sign for different t . For these reasons, we use the fixed basis defined by the time-averaged Hamiltonian for the data presented in Secs. VI. Note that measuring in the fixed basis corresponding to the time-averaged Hamiltonian plausibly mimics “imperfect” quantum dot readout that is slow on the timescale of the typical noise correlation time. In Sec. VII, we circumvent the sign ambiguity problem in the instantaneous basis by noise averaging instead $\langle i\gamma_1(t)\gamma_3(t) \rangle^2$ and are thereby able to directly compare the two measurement approaches (and thus qubit encodings).

V. ANALYTICAL ESTIMATION OF QUBIT TIMESCALES

In this section we use standard techniques to derive analytic formulas for the dephasing and leakage times. These estimates will be compared with timescales extracted from our time-dependent numerical simulations in Sec. VI.

A. Dephasing time

We first restrict attention to the low-energy subspace, neglecting the quasiparticle continuum above the superconducting gap. This approximation is justified provided noise is slow compared to the gap scale. Leakage of the qubit into excited states will be discussed in the next subsection.

Within the low-energy subspace, we model the qubit by the minimal two-level Hamiltonian in Eq. (1). Time dependence in the instantaneous energy splitting $E(t)$ arises due to stochastic temporal variations in parameters in the “parent” microscopic Hamiltonian. We denote fluctuations in these microscopic parameters about their time-averaged values by $\lambda_i(t)$, which are assumed uncorrelated with one another. This set can include variations in the electrochemical potential, magnetic field, etc., in different spatial regions of the device. For example, if only the global chemical potential varies in time by an amount $\delta\mu(t)$, then we simply have $\lambda_1(t) = \delta\mu(t)$; if instead uncorrelated chemical potential variations arise in the left and right halves of the wires, then we have $\lambda_1(t) = \delta\mu_L(t)$ and $\lambda_2(t) = \delta\mu_R(t)$; and so on. Assuming the fluctuations are weak, we Taylor expand the energy to second order in λ_i ’s, yielding

$$E(t) \approx E_0 + \sum_i \lambda_i(t) E'_i + \frac{1}{2} \sum_{i,j} \lambda_i(t) \lambda_j(t) E''_{ij}, \quad (26)$$

where we defined shorthand notation

$$E'_i = \frac{dE}{d\lambda_i} \Big|_{\lambda_i=0}, \quad E''_{ij} = \frac{d^2E}{d\lambda_i d\lambda_j} \Big|_{\lambda_{i,j}=0}. \quad (27)$$

To noise average we will assume that each λ_i exhibits Gaussian noise correlations with mean and variance

$$\overline{\lambda_i(t)} = 0, \quad (28)$$

$$\overline{\lambda_i(t) \lambda_j(t')} = \delta_{ij} S_i(t - t'). \quad (29)$$

We further take $S_i(t)$ to be Gaussian as a function of time, i.e.,

$$S_i(t) = D_i^2 e^{-t^2/(2\tau_i)^2}, \quad (30)$$

corresponding to a power spectrum

$$S_i(\omega) = D_i^2 \sqrt{\frac{4\pi}{\kappa_i^2}} e^{-(\omega/\kappa_i)^2}. \quad (31)$$

In Eqs. (30) and (31), D_i and $\tau_i \equiv 1/\kappa_i$, respectively, denote the fluctuation amplitude and characteristic noise correlation time for fluctuator λ_i ; note that the frequency scale κ_i plays the role of a high-frequency cutoff in the power spectrum. The qubit precession frequency ω_0 then follows from

$$\hbar\omega_0 = \overline{E(t)} = E_0 + \frac{1}{2} \sum_i D_i^2 E''_{ii}. \quad (32)$$

Notice that noise generically shifts the mean energy splitting $\hbar\omega_0$ away from the noise-free value E_0 .

We could in principle also adopt the more realistic case of $1/f$ noise, which leads to similar qualitative conclusions (see, e.g., Ref. [46] for a discussion in a related context). However, in the microscopic models discussed below, $1/f$ noise requires the introduction of a high-frequency cutoff which leads to corrections that make quantitative comparisons between theoretical and numerical estimates challenging.

In the two-level-system approximation used here, it is straightforward to evaluate the readout for the Ramsey-type protocol described in the preceding section. For a particular noise realization one obtains

$$Q(t) \equiv \langle \psi(t) | X_t | \psi(t) \rangle = \cos \left(\int_0^t dt' E(t')/\hbar \right). \quad (33)$$

Appendix B noise averages $Q(t)$; in the long-time limit $t \gg \tau_i$ we find

$$\overline{Q(t \gg \tau_i)} \approx \cos(\omega_0 t) e^{-t/T_2} \quad (34)$$

with dephasing time

$$T_2 = \frac{\hbar^2}{\sqrt{\pi}} \left[\sum_i \tau_i (D_i E'_i)^2 + \frac{1}{2} \sum_{ij} \frac{\tau_i \tau_j}{\sqrt{\tau_i^2 + \tau_j^2}} (D_i E''_{ij} D_j)^2 \right]^{-1}. \quad (35)$$

(Appendix B also examines the short-time limit $t \ll \tau_i$, which is sometimes relevant for our numerical simulations.) Upon varying parameters, e.g., to transition between an ABS and topological qubit, we expect that T_2 will vary by far most dramatically through the E'_i and E''_{ij} factors. These factors quantify the rigidity of the qubit splitting with respect to fluctuations and can be efficiently derived from microscopics through exact diagonalization of *noise-free* static Hamiltonians, in principle containing various bells and whistles. We will carry out such an analysis in Sec. VI in parallel with our explicit time-dependent qubit simulations.

B. Leakage time

We will estimate the qubit leakage time by studying just a single noisy wire in the tetron using Fermi’s golden rule. For analytical tractability, we will further employ a *fixed basis* for Majorana modes in the device; as stressed in Sec. IV B 2, the analytical estimates that follow should thus be viewed as an *upper bound* on the true leakage time that would be obtained

by following the instantaneous basis of maximally localized Majorana modes.

Let us describe the system within the spinful-wire model (extension to the Kitaev-chain description is straightforward). We write the full microscopic Bogoliubov–de Gennes Hamiltonian as $H = H_0 + \delta H(t)$, where H_0 is given by Eq. (5) and $\delta H(t)$ incorporates noise. The static part of the Hamiltonian is diagonalized upon decomposing the fermion $\psi_\sigma(x)$ with spin σ at position x as

$$\psi_\sigma(x) = \Phi_{1\sigma}(x)\gamma_1^{(0)} + \Phi_{2\sigma}(x)\gamma_2^{(0)} + \sum_{E>0} \varphi_{E\sigma}(x)f_E. \quad (36)$$

Here, $\Phi_{1,2\sigma}$ denote the maximally localized Majorana wave functions associated with the fixed-basis Majorana operators $\gamma_{1,2}^{(0)}$ derived from H_0 , while $\varphi_{E\sigma}$ denote wave functions for above-gap excitations created by operators f_E^\dagger . After this decomposition H_0 becomes

$$H_0 = \frac{i}{2}\varepsilon_{12}\gamma_1^{(0)}\gamma_2^{(0)} + \sum_{E>0} E f_E^\dagger f_E \quad (37)$$

with ε_{12} the noise-free splitting of states with $i\gamma_1\gamma_2 = \pm 1$.

We write the noise terms as

$$\delta H(t) = \sum_i \lambda_i(t) \int_0^L dx \psi^\dagger \Lambda_i(x) \psi \quad (38)$$

for some matrices $\Lambda_i(x)$ dependent on the nature of the parameter fluctuations $\lambda_i(t)$. For instance, if $\lambda_1(t) = \delta\mu(t)$ represents a global chemical potential fluctuation, then $\Lambda_1 = -I$. Note that we neglect noise in the pairing channel for simplicity, though such terms could be readily incorporated if desired. Rewriting $\delta H(t)$ using Eq. (36) yields

$$\delta H(t) = \sum_{E>0} \{ [A_{1E}(t)\gamma_1^{(0)} + A_{2E}(t)\gamma_2^{(0)}] f_E + \text{H.c.} \} + \dots \quad (39)$$

The ellipsis denotes pieces involving two Majorana operators or two f_E operators, while the coefficients above read as

$$A_{1,2E}(t) = \sum_i \lambda_i(t) \int_0^L dx \Phi_{1,2}^\dagger(x) \Lambda_i(x) \varphi_E(x). \quad (40)$$

Next, we introduce the complex fermion $f_0 = (\gamma_1 + i\gamma_2)/2$ and define $A_{\pm,E}(t) = A_{1E}(t) \pm iA_{2E}(t)$ so that

$$\delta H(t) = \sum_{E>0} \{ [A_{-,E}(t)f_0 + A_{+,E}(t)f_0^\dagger] f_E + \text{H.c.} \} + \dots \quad (41)$$

According to Fermi's golden rule, the noise-averaged leakage rates Γ_\pm out of the ground states with $f_0^\dagger f_0 = 0$ (Γ_-) and $f_0^\dagger f_0 = 1$ (Γ_+) are given by

$$\Gamma_\pm = \frac{2\pi}{\hbar} \sum_{E>0} \int_{\omega,\omega'} \overline{A_{\pm,E}(\omega)A_{\pm,E}^*(\omega')} \delta(E \mp \varepsilon_{12} - \hbar\omega). \quad (42)$$

The noise correlations in Eqs. (28) through (31) yield

$$\begin{aligned} \overline{A_{\pm,E}(\omega)A_{\pm,E}^*(\omega')} &= 2\pi\delta(\omega - \omega') \\ &\times 2\sqrt{\pi} \sum_j \tau_j D_j^2 e^{-(\omega\tau_j)^2} B_{\pm,j}(E), \end{aligned} \quad (43)$$

where

$$B_{\pm,j}(E) = \left| \int_x [\Phi_1^\dagger(x) \pm i\Phi_2^\dagger(x)] \Lambda_j(x) \varphi_E(x) \right|^2. \quad (44)$$

Feeding this expression into Γ_\pm leads to

$$\Gamma_\pm = 2\sqrt{\pi} \sum_j \frac{\tau_j D_j^2}{\hbar^2} \sum_{E>0} B_{\pm,j}(E) e^{-[(E \mp \varepsilon_{12})\tau_j/\hbar]^2}. \quad (45)$$

Suppose now that E_g is the bulk gap and $\rho(E)$ is the density of states for the above-gap excitations; assuming the coefficients $B_{\pm,j}$ vary smoothly with energy, we approximate the leakage rates as

$$\begin{aligned} \Gamma_\pm &\approx 2\sqrt{\pi} \sum_j \frac{\tau_j D_j^2}{\hbar^2} B_{\pm,j}(E_g) \\ &\times \int_{E_g}^\infty dE \rho(E) e^{-[(E \mp \varepsilon_{12})\tau_j/\hbar]^2}. \end{aligned} \quad (46)$$

Leakage will be most important in the vicinity of the topological phase transition where the bulk gap approaches zero. In this regime, the excitation spectrum can be viewed as arising from gapped-out counterpropagating bulk Majorana fermions with mode velocity v and dispersion

$$E(k) = \sqrt{(\hbar vk)^2 + E_g^2}. \quad (47)$$

The density of states follows as

$$\rho(E) = \frac{L}{\pi \hbar v} \frac{E}{\sqrt{E^2 - E_g^2}} \quad (48)$$

for $E > E_g$ and vanishes otherwise. [We have implicitly assumed periodic boundary conditions in writing Eq. (47), though the choice of boundary conditions should not significantly influence $\rho(E)$.] Assuming $\varepsilon_{12} \ll E_g$, we then obtain

$$\Gamma_\pm \approx \frac{L}{\hbar v} \sum_j \frac{D_j^2}{\hbar} B_{\pm,j}(E_g) e^{-(E_g\tau_j/\hbar)^2}. \quad (49)$$

The coefficients $B_{\pm,j}(E_g)$ depend on details of the wave functions and noise sources. We will use scaling arguments to roughly quantify these factors in three regimes:

(i) *Ultra-short-range noise.* Suppose first that a noise source acts near one of the two Majorana modes, but only on a very local region of size $\xi_j^{\text{noise}} \ll \xi$ (again, ξ denotes the spatial extent of the Majorana wave function). In this case, $\Lambda_j(x)$ in Eq. (44) has appreciable weight only over a distance ξ_j^{noise} so that

$$B_{\pm,j}(E_g) \propto \left(\frac{\xi_j^{\text{noise}}}{\sqrt{\xi L}} \right)^2 = \left(\frac{\xi_j^{\text{noise}}}{\xi} \right)^2 \frac{\xi}{L}. \quad (50)$$

The square root in the denominator simply reflects normalization factors in the Majorana and above-gap-excitation wave functions. Note that $B_{+,j} = B_{-,j}$ here, which implies approx-

imately equal⁶ leakage rates out of the low-energy states with $i\gamma_1\gamma_2 = \pm 1$.

(ii) *Short-range noise.* If $\gamma_{1,2}$ are separated by a sufficiently long distance L_{12} , one can envision a short-range-noise scenario wherein $\xi < \xi_j^{\text{noise}} < L_{12}$. That is, the noise source acts over the entire window of one of the Majorana modes, but does not influence its partner. (This scenario is reasonable for bona fide Majorana zero modes separated by the length of the wire, but is less natural for accidental Andreev bound states.) Here we get

$$B_{\pm,j}(E_g) \propto \left(\frac{\xi}{\sqrt{\xi L}} \right)^2 = \frac{\xi}{L}, \quad (51)$$

again yielding equal leakage rates out of the $i\gamma_1\gamma_2 = \pm 1$ states.

(iii) *Long-range noise.* Finally, consider a long-range-correlated noise source with $\xi_j^{\text{noise}} > L_{12}$. The integral in Eq. (44) now picks up weight from *both* Majorana wave functions, yielding similar scaling to the previous case but with unequal leakage rates for the $i\gamma_1\gamma_2 = \pm 1$ states:

$$B_{\pm,j}(E_g) \propto (1 \pm \eta_j) \frac{\xi}{L} \quad (52)$$

for $-1 < \eta_j < 1$. This regime offers the interesting possibility of a “dark state”: when $\eta_j = \pm 1$ the transition rate for one of the two low-energy states vanishes.

The three cases above can be put on equal footing by writing

$$B_{\pm,j}(E_g) \approx c_j(1 \pm \eta_j) \left(\frac{\xi_j^{\text{noise}}}{\xi} \right)^{b_j} \frac{\xi}{L}, \quad (53)$$

the proportionality constant c_j , exponent b_j , and η_j are determined by properties of noise source j . We thereby obtain leakage times

$$T_{\text{leak}}^{\pm} = 1/\Gamma_{\pm} \approx \hbar^2 \left[\sum_i c_i \frac{\xi}{v} D_i^2 (1 \pm \eta_i) \left(\frac{\xi_i^{\text{noise}}}{\xi} \right)^{b_i} e^{-(E_g \tau_i / \hbar)^2} \right]^{-1}. \quad (54)$$

Comparison to the dephasing time in Eq. (35) (first term) is instructive. We see that the timescale ξ/v appearing in T_{leak}^{\pm} plays the role of τ_i in T_2 ; similarly, the exponential factor $e^{-(E_g \tau_i / \hbar)^2}$ plays the role of the “energetic rigidity” E_i' characterizing the qubit states. Upon approaching the topological phase transition, either from the Andreev-bound-state or true Majorana-zero-mode regime, collapse of the bulk gap E_g readily allows leakage to dominate the qubit lifetime.

VI. NUMERICAL RESULTS

We are now in position to present results on both quasi-analytic evaluation of the dephasing and leakage times (using the formulas derived in Secs. V A and V B) as well as full

microscopic simulations of noisy tetron dynamics for both the Kitaev and spinful nanowire models.

A. Kitaev tetron

We first consider a tetron qubit built from two parallel Kitaev chains [Eq. (4)] each consisting of L physical sites. The time-averaged chemical potentials of the two wires are offset by a small amount μ_{offset} to break degeneracies which would occur for identical wires; that is, $\mu_{\text{top/bot}}^{(0)} = \mu \pm \mu_{\text{offset}}$. We adopt a noise model consisting of two independent fluctuators acting on the respective chemical potentials such that $\mu_{\text{top/bot}}(t) = \mu_{\text{top/bot}}^{(0)} + \delta\mu_{\text{top/bot}}(t)$. Each fluctuator $\delta\mu_{\text{top/bot}}(t)$ obeys Gaussian noise correlations with a Gaussian noise power spectrum as specified in Eqs. (28) through (31). For simplicity, we take the amplitude of typical fluctuations D and noise correlation time $\tau \equiv 1/\kappa$ equal for both fluctuators: $D_{\text{top/bot}} = \delta\mu_{\text{top/bot}}^{\text{typ}} = \delta\mu_{\text{typ}}$ and $\kappa_{\text{top/bot}} = \kappa$.

Given such a setup, it is straightforward to numerically evaluate the analytic (perturbative) estimates for the dephasing and leakage times as derived above in Sec. V. In addition, we perform full microscopic numerical experiments of the noisy tetron dynamics by generating N_{real} independent noise realizations, evolving under the time-dependent Schrödinger equation [which in terms of the covariance matrix is given by Eq. (13)] for each realization, and finally noise averaging the resulting measurements. (Details of our noise generation procedure and subsequent solution of the evolution equation can be found in Appendix A.) As spelled out above in Sec. IV, for our Ramsey-type protocol we initialize the system into a state corresponding to fixed $X_0 = i\gamma_1^{(0)}\gamma_3^{(0)} = +1$, where the four near-zero-energy Majorana modes $\gamma_{1,2,3,4}^{(0)}$ are defined with respect to the time-averaged Hamiltonian A_0 . Throughout the evolution we monitor the expectation values $\langle i\gamma_1\gamma_3 \rangle$, $\langle i\gamma_1\gamma_3 \rangle^2$ (see Sec. IV B 3), and $\langle (i\gamma_1\gamma_2)(i\gamma_3\gamma_4) \rangle$ with $\gamma_{1,2,3,4}$ defined in terms of either the fixed, time-averaged basis used for initialization [i.e., $\gamma_{1,2,3,4} = \gamma_{1,2,3,4}^{(0)}$] or the basis derived from the instantaneous Hamiltonian $A(t)$ [i.e., $\gamma_{1,2,3,4} = \gamma_{1,2,3,4}(t)$]. Finally, we also track the overlap of the evolving wave function onto the low-energy subspace, for both the fixed and instantaneous basis, respectively corresponding to $|a_0|^2 + |b_0|^2$ and $|a_t|^2 + |b_t|^2$ from Eq. (21).

Within a topological phase, both the qubit precession period $2\pi/\omega_0$ and dephasing time T_2 are expected to be exponentially long in the wire length L . Simulating the full microscopic dynamics over these timescales thus becomes prohibitive for modest L even for noninteracting electronic models. Evaluating the relevant timescales using the analytic perturbative predictions, on the other hand, faces no such limitations as this procedure merely requires computing derivatives of the lowest-energy Majorana hybridization energies, which are obtainable with a sparse eigensolver applied to the Hermitian matrix iA .

We present data obtained for numerical experiments performed on a tetron built from two Kitaev wires each containing $L = 10$ physical sites, and with various μ , κ but all other parameters fixed according to $J_{\text{top/bot}} = J = 1$, $\Delta_{\text{top/bot}} = \Delta = 0.4$, $\mu_{\text{offset}} = 0.01$, and, when noise is present, $\delta\mu_{\text{typ}} = 0.02$. (With $\hbar = J = 1$, energies are in units of J , times are in units of \hbar/J , and frequencies, e.g., κ , are in units of J/\hbar .)

⁵We took $\varepsilon_{12}/E_g = 0$ to arrive at Eq. (49); corrections from nonzero ε_{12}/E_g can still give different leakage rates even with short-range-correlated noise.

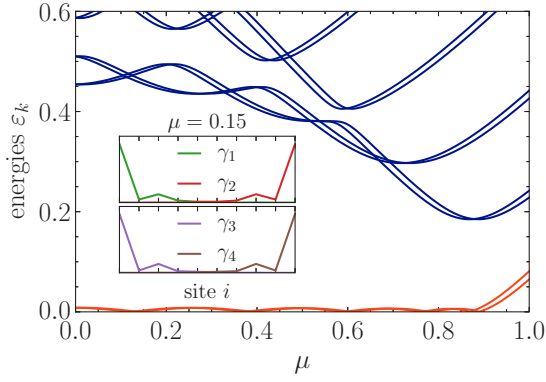


FIG. 3. Energy spectra of the Kitaev tetron as a function of μ for the configuration used in our microscopic numerical simulations (cf. Figs. 4, 5, 10, 11, and 12). Here, each wire of the tetron consists of an $L = 10$ site Kitaev chain with $J_{\text{top/bot}} = J = 1$ and $\Delta_{\text{top/bot}} = \Delta = 0.4$; a slight chemical potential shift of $\mu_{\text{offset}} = 0.01$ is introduced between the two wires to break any associated degeneracies. The low-energy qubit manifold is highlighted in orange, corresponding to the energies $\varepsilon_{1,2}$ in Eq. (9). In the inset, we plot the spatial dependence of the wave-function square modulus for the maximally localized Majorana modes $\gamma_{1,2,3,4}$ (see Sec. IV A) at $\mu = 0.15$.

Figure 3 shows the energy spectra for this qubit configuration as a function of μ in the noise-free limit. Note that for this system size the topological “phase transition” near $\mu = 1$ is significantly smeared by finite-size effects. Within the topological regime at $|\mu| \lesssim 1$, however, fairly well-formed Majorana zero modes arise. The inset plots the spatial profiles of the near-zero-energy maximally localized Majorana modes $\gamma_{1,2,3,4} = \gamma_{1,2,3,4}^{(0)}$ at $\mu = 0.15$.

Figure 4 contains simulation results for a noisy system with $\mu = 0.15$ and $\kappa = 0.05$. The top panel shows the evolution of $\langle i\gamma_1\gamma_3 \rangle$ —defined here in terms of the fixed basis shown in the inset of Fig. 3—for 10 independent noise realizations. Also shown for comparison are data for a noiseless run (thick black curve). The bottom panel shows the results of averaging $N_{\text{real}} = O(10^3)$ noise realizations,⁷ a fit of this noise-averaged numerical data to the functional form $\cos(\omega_0 t)e^{-t/T_2}$, as well as the time dependence predicted⁸ by the (small-noise-amplitude) analytic calculations of Sec. V A [cf. Eqs. (34) and (35)]. We see excellent agreement between the numerical data and analytical prediction. These parameters yield negligible leakage (as determined by the relevant measurements discussed above; not shown) on the timescale of the dephasing time. That is, $T_{\text{leak}} \gg T_2$.

Finally, in Fig. 5, we present results of a comprehensive study varying μ at different values of $\kappa = 0.001, 0.01, 0.05$, and 0.1 . The bottom panels plot the qubit precession period $2\pi/\omega_0$ versus μ as obtained by both the microscopic

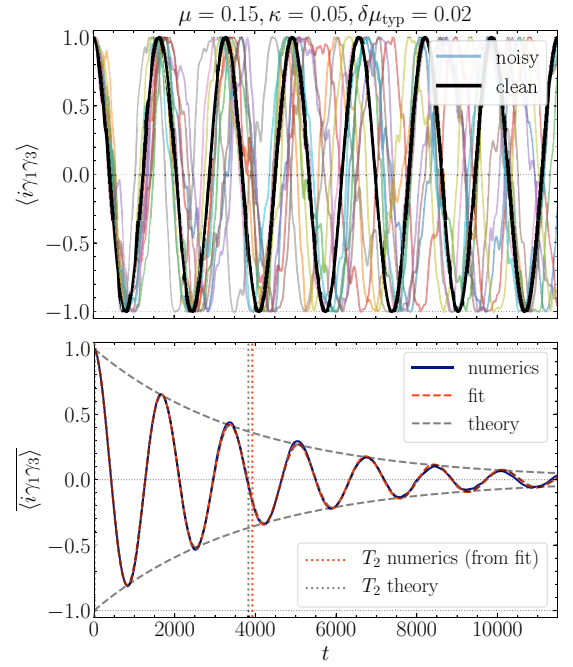


FIG. 4. Example single noise realizations (top) and noise-averaged data (bottom) for the (quantum-averaged) quantity $\langle i\gamma_1\gamma_3 \rangle$ (with $\gamma_{1,3} = \gamma_{1,3}^{(0)}$) simulated in the Kitaev tetron of Fig. 3, here at $\mu = 0.15$. The top panel shows the results of 10 independent noise realizations (“noisy”; light curves) with $\kappa = 0.05$ and $\delta\mu_{\text{typ}} = 0.02$ (one global μ fluctuator on each wire of the tetron), as well as for a system without noise (“clean”; black curve). In the bottom panel, we present the noise-averaged signal (“numerics”; solid blue) and a fit (“fit”; dashed orange) guided by the theoretical prediction of Eq. (34). The theoretical prediction itself (and its envelope) obtained using Eqs. (32)–(35) is also plotted (“theory”; dashed gray). The orange and gray vertical dotted lines indicate the respective T_2 times.

numerical simulations (“numerics”) and analytic prediction (“theory”) [see Eq. (32) and note that the predicted qubit precession frequency is independent of κ]. The top panels similarly show the corresponding coherence times T_2 and, in the rightmost plot, T_{leak} . Numerical T_2 data correspond to fits of the noise-averaged, fixed-basis $\langle i\gamma_1\gamma_3 \rangle$ data to an oscillatory exponential, precisely as in Fig. 4. For the analytic T_2 predictions, we show the timescales predicted both in the limit $t \gg 1/\kappa$ (“ T_2 theory”) and $t \ll 1/\kappa$ [“ T_2 theory (transient)”]. In the transient case, the decay is expected to follow a Gaussian instead of an exponential [see Eqs. (B16) and (B17) in Appendix B]. As a guide, in the left two panels where the noise is “slow,” the horizontal green lines indicate the timescale $\tau = 1/\kappa$. Coherence times falling below that scale indicate importance of the transient regime. For consistency, however, we fit the data for all numerical experiments to an exponential form. Nonetheless, when the coherence times fall below $\tau = 1/\kappa$, the extracted timescales from the numerics track the $t \ll 1/\kappa$ prediction reasonably well. To summarize, when the dephasing time measured in the numerics is much larger than $1/\kappa$, it matches the (long-time) analytic prediction very well; and when it dips below $1/\kappa$, it follows the (short-time, “transient”) analytic prediction.

⁷Error bars are typically on the order of the symbol size or smaller for all data that we present.

⁸Throughout, when evaluating the energy splitting E in Eqs. (35) and (32) for a given microscopic model, for simplicity we take $E = \varepsilon_1 + \varepsilon_2$ [with $\varepsilon_{1,2} \geq 0$, cf. Eq. (9)] without enforcing a fixed global parity.

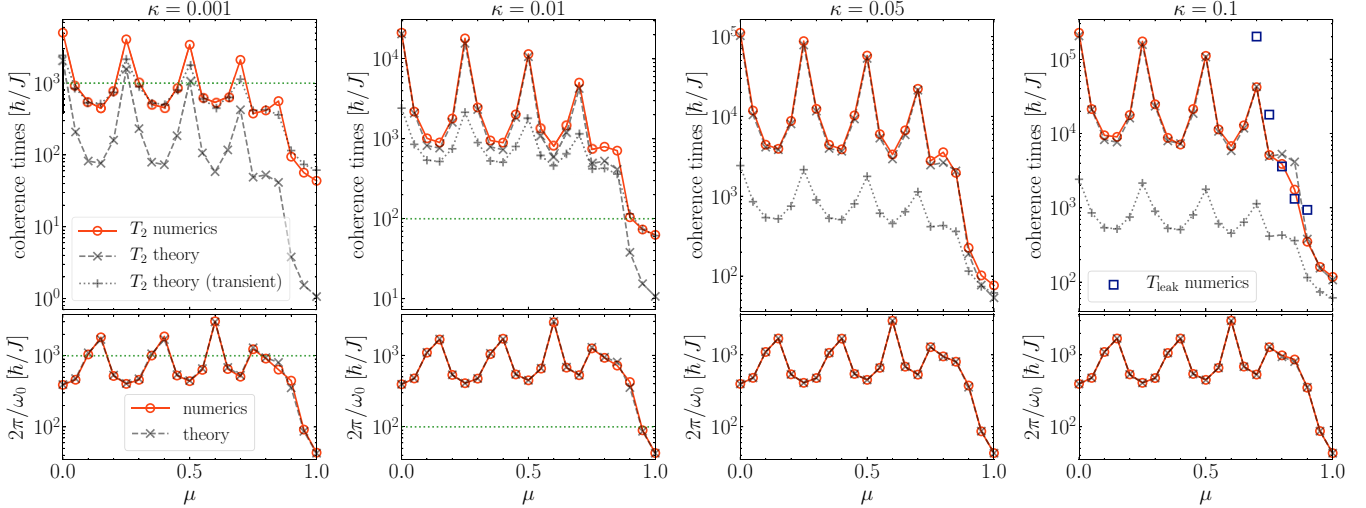


FIG. 5. Numerically extracted coherence times (top panels) and qubit precession periods (bottom panels) from full microscopic numerical simulations of a noisy Kitaev tetron (see Figs. 3 and 4 for details of the setup). For the dephasing timescales, we plot in the top panels both the values obtained numerically via measurement of $\langle i\gamma_1\gamma_3 \rangle$ (“ T_2 numerics”) as well as the analytic predictions of Eqs. (35) and (B17) [“ T_2 theory” and “ T_2 theory (transient),” respectively]; similarly, in the bottom panels, we show both numerically extracted (“numerics”) and analytically predicted [Eq. (32)] (“theory”) precession periods. From left to right, the four panels correspond to increasing $\kappa = 0.001, 0.01, 0.05, 0.1$ (roughly the high-frequency cutoff of the noise power spectrum) or, equivalently, decreasing noise correlation times $\tau = 1/\kappa$ (the dotted green lines in the two leftmost panels indicate the timescale τ). For all data, we take independent global chemical potential fluctuators on each wire of the tetron, each with typical fluctuation amplitude $\delta\mu_{\text{typ}} = 0.02$. Only at $\kappa = 0.1$ and near the phase transition out of the topological phase for $\mu \sim 1$ does the system exhibit detectable leakage times (“ T_{leak} numerics”); see text for more details.

As in the example presented in Fig. 4, for all of the data points with $\kappa \leq 0.05$ in Fig. 5, we detect negligible leakage out of the low-energy qubit subspace over the timescale of dephasing. Only for $\kappa = 0.1$ at $\mu \sim 1$ upon exiting the topological phase do we observe appreciable leakage. The quantities T_{leak} in the rightmost panel of Fig. 5 were extracted by fitting $|a_t|^2 + |b_t|^2$ [see Eqs. (21) and (22)] to the (phenomenological) form $Ae^{-t/T_{\text{leak}}} + (1 - A)$; only data points for which $A > 0.1$ are plotted. We see from the energy spectra in Fig. 3 that the dip in leakage times for $\mu \lesssim 1$ coincides with the finite-size crossover into the trivial phase on this small $L = 10$ site Kitaev tetron, i.e., the point where the gap is minimal. On the other hand, for small $\kappa \leq 0.05$ the minimum finite-size gap exceeds κ so that the qubit lifetime is governed by the dephasing time T_2 for all μ .

B. Spinful nanowire tetron

We now turn to tetrons assembled from two spinful nanowires governed by the Hamiltonian defined in Eq. (5). Guided by the data presented in Ref. [25], we focus on a set of parameters that yields low-energy edge ABSs over a sizable window of Zeeman fields $V'_{zc} < V_z < V_{zc}$ before entering a topological regime with bona fide Majorana zero modes at larger fields $V_z > V_{zc}$. Above in Fig. 2, we presented data of qubit coherence times and precession periods for such a system as a function of V_z evaluated numerically according to the analytical estimates of Sec. V. The corresponding (time-averaged) system parameters are as follows: $\Delta(x) = \frac{\Delta_0}{2} [\tanh(\frac{x-x_0}{\ell_\Delta}) + 1]$ with $\Delta_0 = 0.25$ meV, $\ell_\Delta = 0.03$ μm , and $x_0 = 0.3$ μm ; $V(x) = \frac{V_0}{2} [-\tanh(\frac{x-x_0}{\ell_V}) + 1]$ with $V_0 = 3.8\Delta_0$ and $\ell_V = 0.03$ μm ; $m = 0.03m_e$; $\mu = V_0 = 3.8\Delta_0$

(with top/bottom wire offset $\mu_{\text{offset}} = 0.01$ meV); and $\alpha = 500$ meV \AA . For our discretization, we take a lattice spacing $a \approx 0.01$ μm corresponding to $N_{\text{sites}} = 100, 200, 300, 400$ sites (per wire) for the sequence of sizes $L = 1, 2, 3, 4$ μm . In Fig. 6, we show additional data for the $L = 2$ μm system. The top panel plots the energy spectrum versus V_z while the bottom panels illustrate the spatial profiles of the maximally localized near-zero-energy Majorana modes $\gamma_{1,2,3,4}$ at $V_z = 0.6$ meV (nontopological ABS qubit) and 1.4 meV (topological qubit). The chosen noise configuration consists of four independent fluctuators: one global, spatially constant chemical potential (μ) and Zeeman field (V_z) fluctuator per wire each with Gaussian correlated noise with inverse correlation times $\kappa = 0.05$ meV/ \hbar and typical fluctuation amplitudes $\delta\mu_{\text{typ}} = \delta V_{z,\text{typ}} = 0.01$ meV.

Given these parameters, the T_2 and ω_0 values plotted in Fig. 2 were obtained by numerically evaluating Eqs. (35) and (32) via sparse diagonalization of the Hermitian matrix iA and computing the requisite energy derivatives. (Only the leading-order terms were included in the evaluations for the more conceptual Fig. 2, while for Figs. 5 and 8 the second-order terms were also calculated and are in fact necessary to obtain quantitative agreement with the microscopic numerics.) For T_{leak} we take a simplified approach which focuses entirely on the predicted exponential dependence: $T_{\text{leak}} \sim e^{(E_g/\hbar\kappa)^2}$. Specifically, we evaluate Eq. (54) at fixed $c_i = 1$, $\eta_i = 0$, $b_i = 0$, $v = \alpha/\hbar (= 500$ meV $\text{\AA}/\hbar)$, and $\xi = 200$ nm (recall $D_i = \delta\mu_{\text{typ}} = \delta V_{z,\text{typ}} = 0.01$ meV and $\kappa_i = \kappa = 0.05$ meV/ \hbar). While the analysis of Sec. VB considered only a single nanowire, it can be safely applied to the noisy tetron considered here as the wires are taken to be decoupled and the spatial profiles of all noise sources have

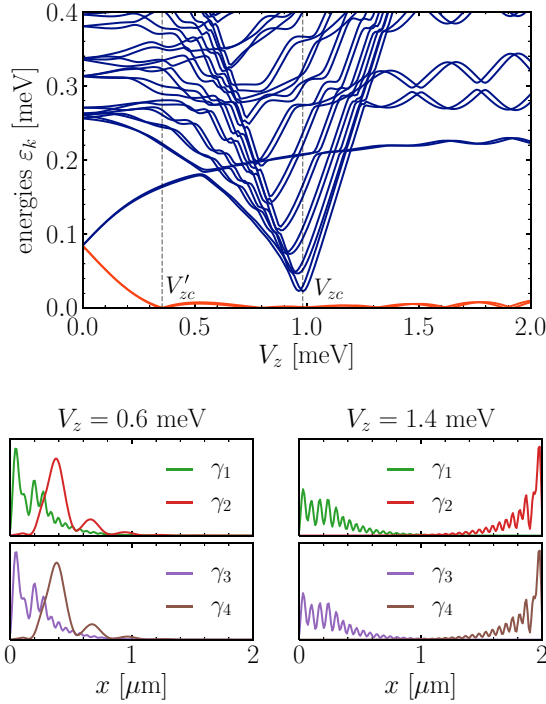


FIG. 6. Energy spectra (top) and representative near-zero-energy Majorana wave functions (bottom) for the continuum-limit $L = 2 \mu\text{m}$ spinful tetron configuration analyzed in Fig. 2. The chosen parameters were inspired by Ref. [25] (see text for all details) to give a sizable window of a nontopological ABS regime for $V'_{zc} < V_z < V_{zc}$ before the onset of the topological phase transition at V_{zc} . In the bottom panels, we show the square modulus of the wave functions for the maximally localized Majorana modes $\gamma_{1,2,3,4}$ (summing both spin components) at $V_z = 0.6$ meV (partially separated ABS qubit) and $V_z = 1.4$ meV (topological qubit).

support on only one wire or the other (two global fluctuators per wire). In evaluating Eq. (54), we thus consider only two fluctuators, and for the excitation gap E_g , we take an average of the excitation gaps $[\varepsilon_{3,4}]$ from Eq. (9) for the top and bottom wires (recall the finite μ_{offset}). This procedure ignores the complicated dependencies of the (qualitatively less important) subexponential factors on the wave-function amplitudes, noise profile details, etc., but it faithfully captures the dominant exponential dependence and thus suffices for our purposes here.

As highlighted in Sec. II, Fig. 2 demonstrates several important points: First, within the ABS regime ($V'_{zc} < V_z < V_{zc}$) the qubit coherence time and precession frequency exhibit weak, nonuniversal dependence on wire length L . Second, the qubit lifetime drops sharply near the topological phase transition ($V_z \sim V_{zc}$) due to noise-induced leakage out of the qubit subspace, leading to a leakage-limited qubit. Third, deep in the topological phase the system forms an exponentially protected, dephasing-limited topological qubit (though for still longer wires the topological qubit will eventually become leakage limited as the bulk excitation gap is roughly independent of L). Fourth, within the topological phase, the dephasing time and qubit precession period oscillate out of phase and eventually decrease upon increasing V_z , reflecting an enlargement of the coherence length ξ at “large” V_z .

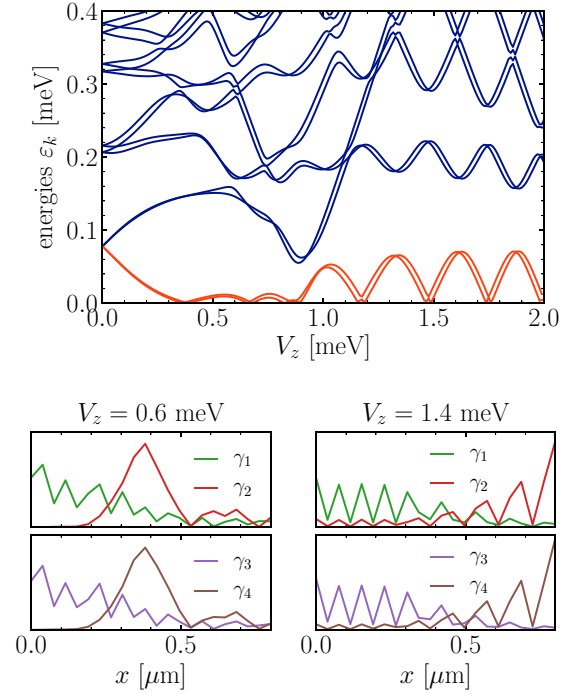


FIG. 7. Energy spectra (top) and representative near-zero-energy Majorana wave functions (bottom) for the $L = 0.8 \mu\text{m}$ spinful tetron configuration studied in Fig. 8; these data are analogous to Fig. 6 but with some parameter changes (see text for all details). Here we discretize the spinful nanowire model of Eq. (5) onto $N_{\text{site}} = 22$ sites per wire to make subsequent full microscopic simulations of the noisy dynamics numerically tractable, while still maintaining several of the qualitative features of the continuum-limit model of Figs. 2 and 6.

Performing full microscopic simulations of such noisy spinful tetrons containing hundreds of lattice sites in a manner that parallels the Kitaev simulations presented in Fig. 5 is numerically intractable. We can, however, construct a small- L “toy” system far from the continuum limit of Eq. (5) that is amenable to simulations and exhibits qualitatively similar features to the systems considered in Fig. 2. For this system, we choose a set of parameters [including functional forms of the external potentials $V(x)$ and $\Delta(x)$] identical to those used in Figs. 2 and 6 with the following exceptions: now $\alpha = 400 \text{ meV } \text{\AA}$, $\ell_V = 0.08 \mu\text{m}$, and $L = 0.8 \mu\text{m}$ discretized into $N_{\text{sites}} = 22$ (per wire) such that the “lattice constant” $a = \frac{L}{N_{\text{sites}} - 1} \approx 0.038 \mu\text{m}$. In Fig. 7, we show the corresponding energy spectra as a function of external Zeeman field V_z as well as the wave-function amplitudes for the maximally localized Majorana modes $\gamma_{1,2,3,4}$ at $V_z = 0.6$ meV (ABS qubit) and 1.4 meV (topological qubit). For the noise setup, we again take one global μ and V_z fluctuator on each wire of the tetron with typical fluctuation amplitudes $\delta\mu_{\text{typ}} = \delta V_{z,\text{typ}} = 0.01 \text{ meV}$ and identical inverse correlation times κ .

Using these parameters, we show in Fig. 8 data for a spinful tetron analogous to the Kitaev tetron results of Fig. 5. Here, as in Fig. 2, we sweep over V_z ; the four panels correspond to $\kappa = 0.001, 0.01, 0.05, 0.1 \text{ meV}/\hbar$. The extraction of the numerical (fitted) timescales (numerics) and evaluation of the analytical timescales (theory) were carried out in exactly

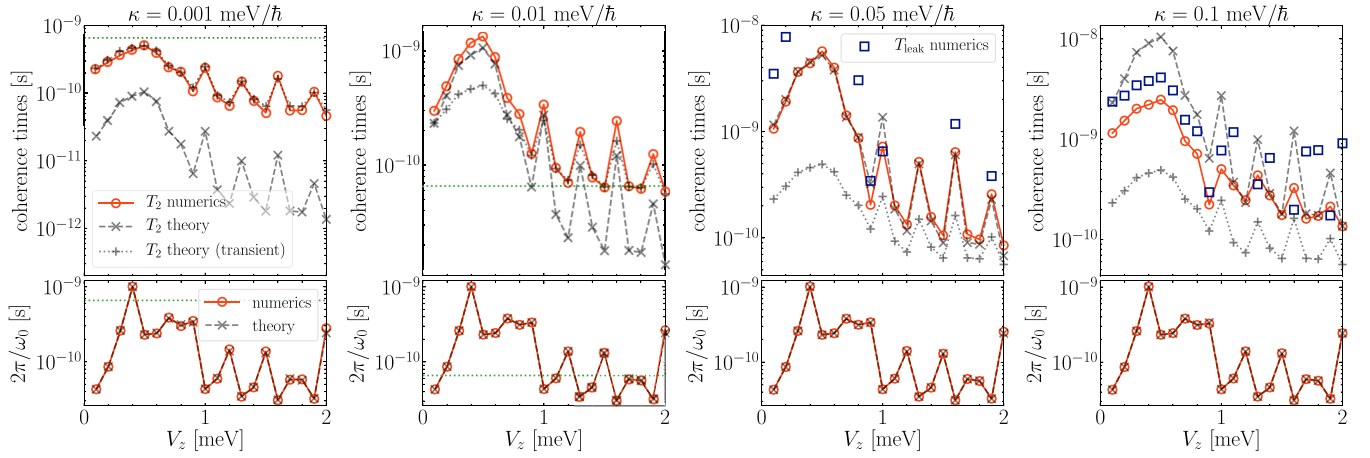


FIG. 8. Numerically extracted coherence times (top panels) and qubit precession periods (bottom panels) from full microscopic numerical simulations of a noisy spinful nanowire tetron (see text and Fig. 7 for details of the setup). These data are analogous to the Kitaev tetron data of Fig. 3; and all formatting, conventions, and means of data analysis are identical. Again, the presence of numerically obtained T_{leak} points at $\kappa = 0.05, 0.1$ meV/ \hbar indicate a detectable leakage time via measurement of $|a_r|^2 + |b_l|^2$. Importantly, when this timescale drops below the analytically predicted T_2 time, the numerically extracted T_2 time (obtained via measurement of $\langle i\gamma_1\gamma_3 \rangle$) itself gets cut off and drops below the analytic value.

the same manner as in Fig. 5. Again, in the two leftmost panels at small $\kappa = 0.001, 0.01$ meV/ \hbar we (1) see the importance of the transient regime when comparing numerical T_2 times to the theoretical predictions and (2) observe no detectable leakage out of the ground-state manifold on the simulated timescales for all V_z . As seen in Fig. 8, the crossover from the ABS regime to topological regime occurs near $V_z \approx 0.9$ meV where the excitation gap is minimal. In this vicinity, for $\kappa = 0.05$ meV/ \hbar , we see that the numerically obtained leakage times become on the order of or smaller than the theoretically predicted dephasing times. In turn, the *numerical* dephasing times drop below the analytic prediction: the noise-averaged signal $\langle i\gamma_1\gamma_3 \rangle$ is now damped by both dephasing effects (cf. Fig. 4) as well as leakage effects in which the amplitude of oscillations for *individual noise realizations* also decay in time (see also Fig. 11). While our theoretical analysis in Sec. V treated the effects of dephasing and leakage independently, our full microscopic simulations as presented here are able to capture both effects simultaneously and faithfully. (The same physics is also occurring near the crossover out of the topological phase in the rightmost panel of Fig. 5.)

Finally, at $\kappa = 0.1$ meV/ \hbar in the rightmost panel of Fig. 8, we observe an intricate interplay between dephasing and leakage. Here, κ is sufficiently large to cutoff the numerically extracted T_2 time below the pure dephasing limit even in the ABS regime for $V_z \lesssim 0.9$ meV. In the vicinity of the crossover we see a sharp reduction in the leakage time due to gap closing, thereby limiting the qubit lifetime as discussed above in the context of Fig. 2. (Note that for the particular qubit simulated in Fig. 8, due to, e.g., the extremely short wire length, even pure dephasing effects cause a reduction in coherence times in the topological regime relative to the ABS one; there is thus a concomitant drop near the corresponding crossover which is unrelated to the leakage-induced qubit lifetime reduction due to gap closing.) Within the topologi-

cal phase, for points corresponding to a minimal excitation gap [for example, at $V_z = 1.3, 1.6, 1.9$ meV (cf. Fig. 6)], $\kappa = 0.1$ meV/ \hbar is again sufficiently large to cause leakage to largely dictate the qubit coherence time. (In principle, leakage-limited lifetimes deep within the topological phase could also occur for sufficiently fast noise in the continuum limit analytical study of Fig. 2; however, studying these effects quantitatively in that framework would require a less crude evaluation of the analytical T_{leak} prediction.) On the other hand, for other values of V_z with excitation gaps in excess of $\hbar\kappa = 0.1$ meV, the qubit lifetime is limited by dephasing, and the numerically extracted timescales match the analytic T_2 prediction impressively well.

C. Estimation of coherence times for realistic nanowire models

A key advantage of the approach outlined in Sec. VA to estimate dephasing times is that it relies purely on the derivative of the Majorana splitting with respect to the fluctuating variables. Such simple spectral properties can be obtained using sparse diagonalization techniques and are thus easy to evaluate also for models where a computation of the full dynamics would be prohibitively expensive. To further illustrate this point, we perform such an estimate for a more sophisticated model of a nanowire that includes a realistic electrostatic potential and several subbands. We closely follow the approach of Ref. [71], which considers a rectangular wire of dimensions $L_x \times L_y \times L_z$ proximitized on the top by a superconductor of thickness d_z ; here, z is the direction transverse to the superconductor and x is the long direction of the wire, i.e., $L_x \gg L_y, L_z$. We simplify the model in two key ways: First, we ignore the effect of subbands in the y direction. And second, instead of treating the superconductor explicitly, we directly induce superconductivity in the semiconductor through a mean-field pairing term. We perform simulations with fixed $L_z = 60$ nm and variable L_x .

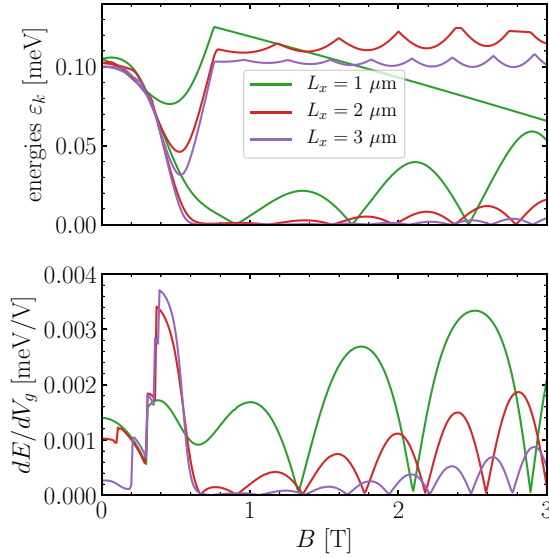


FIG. 9. Dephasing analysis of “slab geometry” nanowire. The upper panel shows the low-energy spectrum for a single wire, while the lower panel shows the derivative of the qubit splitting E with respect to gate voltage V_g . Given some knowledge about the noise environment, this quantity can be directly related to the dephasing time T_2 via Eq. (35). For a detailed list of parameters used in these simulations, see main text. We find that also for this more realistic, multiband model of a nanowire, the qualitative features such as exponential dependence of dephasing time on the length of the wire as well as the correlation between splitting amplitude and dephasing time are preserved.

The first-quantized normal-state Hamiltonian of the system can be written as

$$H_N = -\frac{1}{2m^*}(\partial_x^2 + \partial_z^2) - \alpha \hat{k}_x \sigma_y + \phi(z) + \frac{\mu_B g B}{2} \sigma_x, \quad (55)$$

with m^* the effective mass, α the strength of spin-orbit coupling, $\phi(z)$ the electrostatic potential, g the g factor of the semiconducting material, B the external magnetic field, and σ_α the Pauli matrices acting on spin. We use $m^* = 0.026$, $g = -15$, $\alpha = 0.05$ eV nm. Using a Nambu-space notation, where τ_α are Pauli matrices acting in particle-hole space, we can write the Hamiltonian of the superconducting system as

$$H = H_N \tau_z - \Delta \sigma_y \otimes \tau_y, \quad (56)$$

where Δ denotes the strength of superconducting pairing, which we set to $\Delta = 0.1$ meV. The electrostatic potential $\phi(z)$ is obtained from a self-consistent Thomas-Fermi calculation, where the boundary condition near the superconductor is set to an assumed band offset of 300 meV, while the boundary condition at the other end is tuned via an electrostatic gate with an applied voltage V_g . We use a gate voltage of $V_g = -0.288$ V, which tunes the system close to the bottom of a band and thus favors the formation of a topological phase.

The spectrum is evaluated by employing a finite-difference approximation, with a regular discretization of $a = 2$ nm, and using a shift-and-invert eigensolver. Results are shown in Fig. 9. Here, we evaluate the derivative of the energy splitting

E with respect to the applied gate voltage V_g . Using Eq. (35), this estimate can be combined with an estimate for the noise correlation time as well as typical fluctuation amplitude to obtain an estimate for the T_2 dephasing time. We note that this calculation can in principle be extended to more realistic models of the system, including explicit treatment of the superconductor and even full three-dimensional models of the device. However, major uncertainty enters the estimates through the properties characterizing the noise, which can only be extracted from experiments.

VII. EFFECTS OF FIXED VERSUS INSTANTANEOUS MEASUREMENT BASIS

In this section, we investigate in more quantitative detail the effects of encoding and measuring the qubit in the basis of maximally localized Majorana modes associated with the *instantaneous* Hamiltonian, as opposed to using a *fixed* basis throughout the time evolution. In our numerical simulations presented in Sec. VI, we used the fixed basis of maximally localized Majorana modes $\gamma_{1,2,3,4} = \gamma_{1,2,3,4}^{(0)}$ corresponding to the *time-averaged* Hamiltonian specified by A_0 to initialize the qubit. Furthermore, for technical reasons described in Sec. IV B 3, we often employed this same basis for readout of quantities such as $\langle i\gamma_1 \gamma_3 \rangle$. For a given noise realization, the initial Hamiltonian will differ from the time-averaged one, i.e., $A(t=0) \neq A_0$, and hence this procedure introduces a sort of “quench” in the dynamics at $t=0$, the effects of which become more pronounced for large typical fluctuation amplitudes D_i . For the small, seemingly reasonable D_i chosen above, this effect is negligible. We expect, however, the difference in measurement basis choice to manifest itself most clearly when D_i becomes large. Here, it is worth reiterating an important point made in Sec. IV B 2: Fluctuations, especially those of large amplitude, may significantly alter the maximally localized Majorana modes (e.g., their locations), but in the purely adiabatic limit the qubit subspace is still preserved provided that one tracks and measures in the instantaneous basis. On the other hand, measuring in a fixed basis throughout the evolution will naturally underestimate qubit coherence times, as the noisy dynamics will generically induce a “leakage” out of the fixed basis over time, even if the qubit subspace is perfectly preserved by the instantaneous basis.

Initializing using the time-averaged basis $\gamma_{1,2,3,4}^{(0)}$ is clearly problematic for addressing these matters at large D_i ; e.g., even at $t=0$ the system may have measurable leakage out of the qubit subspace derived from the instantaneous basis. For this section, we thus employ a different initialization procedure from that used in Sec. VI: For each noise realization, we find a maximally localized set of near-zero-energy Majorana modes $\gamma_{1,2,3,4}(t=0)$ derived from $A(t=0)$; we subsequently set $i\gamma_1(t=0)\gamma_3(t=0) = +1$ and $i\gamma_2(t=0)\gamma_4(t=0) = \pm 1$, with the latter chosen such that the global fermion parity is always fixed to $+1$. We then compare the results of two different measurement bases: (1) the initial basis defined by the $\gamma_{1,2,3,4}(t=0)$ (which is fixed for each noise realization) and (2) the instantaneous basis defined by $\gamma_{1,2,3,4}(t)$. [For small enough D_i , we have checked that using the former initial/fixed measurement basis produces results which are

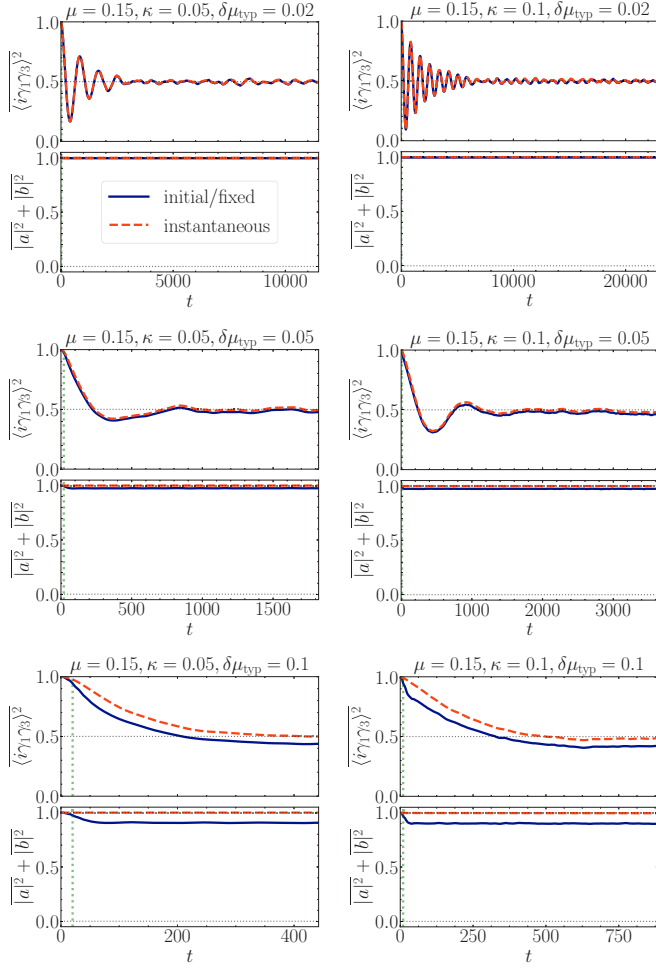


FIG. 10. Measurements of $\overline{\langle i\gamma_1\gamma_3 \rangle^2}$ and $\overline{|a|^2 + |b|^2}$ in both the fixed/initial (solid blue) and instantaneous (dashed orange) basis for a Kitaev tetron at $\mu = 0.15$, $\kappa = 0.05, 0.1$ (increasing left to right), and $\delta\mu_{\text{typ}} = 0.02, 0.05, 0.1$ (increasing top to bottom); see text for details of other chosen parameters. Here, we initialize the system using a procedure based on the *initial* Hamiltonian encoded by $A(t=0)$. For large enough $\delta\mu_{\text{typ}}$, a noticeable difference in the choice of measurement basis is observed on a timescale on the order of $\tau = 1/\kappa$ (vertical green dotted lines).

numerically indistinguishable from the initialization/readout procedure used in Sec. VI based on the $\gamma_{1,2,3,4}^{(0)}$.

Figures 10 and 11 present examples of noise-averaged time dynamics of a Kitaev tetron comparing the results of the two different measurement bases. We focus on the following two quantities: $\overline{\langle i\gamma_1\gamma_3 \rangle^2}$ and $\overline{|a|^2 + |b|^2}$, where for the initial/fixed measurement basis $\gamma_i = \gamma_i(t=0)$ and $|a|^2 + |b|^2 = |a_{t=0}|^2 + |b_{t=0}|^2$, while for the instantaneous basis $\gamma_i = \gamma_i(t)$ and $|a|^2 + |b|^2 = |a_t|^2 + |b_t|^2$. (We consider $\overline{\langle i\gamma_1\gamma_3 \rangle^2}$ instead of $\overline{\langle i\gamma_1\gamma_3 \rangle}$ as the latter would be polluted by the sign ambiguity described in Sec. IV B 3 in the case of the instantaneous basis; see also Fig. 12.) The noise model is the same as that considered in Sec. VIA: one global μ fluctuator on each wire of the tetron (for a total of two fluctuators), each characterized by typical amplitude $\delta\mu_{\text{typ}}$ and inverse correlation time κ . We perform simulations with $\delta\mu_{\text{typ}} = 0.02, 0.05, 0.1$ and $\kappa = 0.05, 0.1$. Figure 10 corresponds to $\mu = 0.15$ which is deep in the

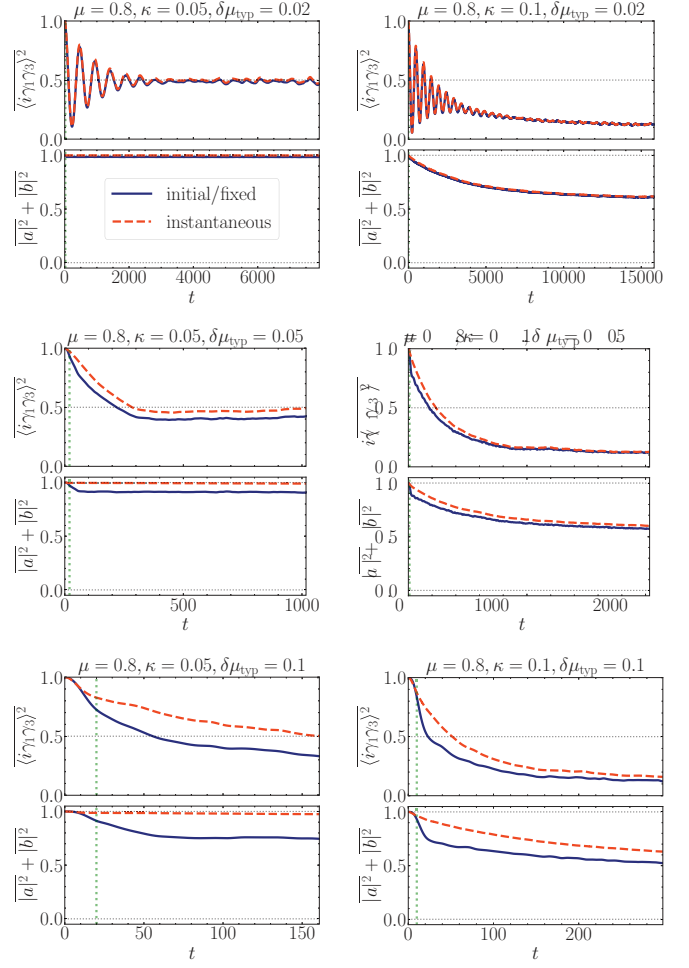


FIG. 11. Data analogous to that presented in Fig. 10 but here taken at $\mu = 0.8$. Now leakage out of the instantaneous low-energy qubit subspace is observed at $\kappa = 0.1$ (as detected by $|a_t|^2 + |b_t|^2$). For all other data at $\mu = 0.15$ in Fig. 10 and $\kappa = 0.05$ here, leakage out of the instantaneous qubit subspace is negligible. Measuring in the fixed basis, on the other hand, generically leads to an effective (topologically unprotected) leakage out of the fixed-basis manifold, which can become significant for large $\delta\mu_{\text{typ}}$, thereby underestimating qubit coherence times.

topological phase, while Fig. 11 corresponds to $\mu = 0.8$ near the crossover into the trivial phase (see Fig. 3). The top panels show the time dependence of $\overline{\langle i\gamma_1\gamma_3 \rangle^2}$, while the bottom panels show $\overline{|a|^2 + |b|^2}$; solid blue (dashed orange) curves represent measurements in the initial/fixed (instantaneous) basis. In the case of pure dephasing, $\overline{\langle i\gamma_1\gamma_3 \rangle^2}$ is expected to take the following form (for times $t \gg 1/\kappa$): $\frac{1}{2}[\cos(2\omega_0 t)e^{-4t/T_2} + 1]$, i.e., relative to $\overline{\langle i\gamma_1\gamma_3 \rangle}$ [cf. Eq. (34)] it oscillates at twice the frequency and approaches $\frac{1}{2}$ (instead of 0) at four times the decay rate.

For relatively small $\delta\mu_{\text{typ}} = 0.02$, we see in the first row of data presented in Figs. 10 and 11 that the chosen measurement basis negligibly impacts the results.⁹ This conclusion holds

⁹The residual oscillations in $\overline{\langle i\gamma_1\gamma_3 \rangle^2}$ at long times are a convergence artifact associated with noise averaging the square.

both in the case of (nearly) pure dephasing dynamics ($\mu = 0.15$ at $\kappa = 0.05, 0.1$ and $\mu = 0.8$ at $\kappa = 0.05$) and dynamics involving both dephasing and leakage ($\mu = 0.8$ at $\kappa = 0.1$). Note that in the latter case, $\langle i\gamma_1\gamma_3 \rangle^2$ approaches a value less than $\frac{1}{2}$ as the oscillations die out: leakage causes the traces for individual noise realizations to themselves decay in oscillation amplitude over time (in contrast to Fig. 4).

As we increase $\delta\mu_{\text{typ}}$ in the middle ($\delta\mu_{\text{typ}} = 0.05$) and bottom ($\delta\mu_{\text{typ}} = 0.1$) rows of Figs. 10 and 11, we see that the difference in measurement basis becomes more pronounced. In addition, at fixed $\delta\mu_{\text{typ}}$ this difference is most severe for slow noise; this point is particularly apparent in the bottom row of Fig. 11 comparing $\kappa = 0.05$ to 0.1. Furthermore, for both considered measurements, we note that the difference becomes manifest only after a timescale on the order of the noise correlation time $\tau = 1/\kappa$ (indicated by vertical green dotted lines). Ultimately, we can clearly see from this data that the relevant coherence time diagnostics decay more quickly when evaluated in a fixed basis.¹⁰ These results are fully consistent with the qualitative picture sketched above in Sec. IV.

VIII. DISCUSSION

In this paper we explored a Ramsey-type protocol that probes qubit dynamics in proximitized nanowire devices, which can support either a topological Majorana-based qubit or a trivial ABS qubit depending on parameters. These two scenarios are challenging to distinguish in local probes, e.g., transport, yet display vastly different noise sensitivity as quantified by the qubit dephasing time revealed by our protocol. The required measurements are more challenging than transport but yield correspondingly more detailed information including the qubit lifetime, time-domain detection of the qubit splitting, and the presence of a topological phase transition. In our study we employed both analytical estimates of dephasing and leakage times as well as explicit numerical simulations of the noisy qubit dynamics. An appealing feature of the analytical estimates is that they can be readily evaluated (modulo uncertainties in noise details that require experimental input) even in state-of-the-art microscopic models. Some proof-of-concept simulations were presented in Sec. VIC, and it would be valuable to further quantify the “rigidity” of the qubit splitting to fluctuations in future modeling efforts. Another feature highlighted by our study is the distinction between fixed-basis and instantaneous Majorana modes in a noisy environment. We argued on general grounds, and confirmed in our explicit simulations, that examining the former underestimates the true qubit lifetime, with the effect becoming increasingly prominent as the noise becomes slower and of larger amplitude.

In actual experiments, additional imperfections could obfuscate some of the features of the dephasing time discussed in this paper. For example, due to long-range inhomogeneities (on a scale longer than the superconducting coherence length) in the electrostatic potential and other system parameters,

a nanowire may not undergo a topological phase transition simultaneously at all positions. Instead, the critical field may vary in space, so that some regions can enter the topological phase earlier than others. In such cases, the minimal gap when sweeping the magnetic field does not scale inversely with the system size, but instead inversely with the size of the largest contiguous region undergoing the phase transition. However, the analysis of Sec. VB shows that the relevant regions would be those whose critical states significantly overlap with the Majorana wave functions, i.e., regions proximate to the ends of the wire. Thus, the leakage time near the phase crossover regime will be determined by the properties of those critical regions.

Moreover, throughout this paper (e.g., in Fig. 2) we have ignored magnetic-field-induced suppression of the bulk superconducting gap. Such effects are clearly important in present-day experiments [10,11] and may ultimately limit the feasibility of our proposed fixed-length study in those devices. Nevertheless, we are hopeful that future devices will harbor a larger window of field strengths over which one can scrutinize trends in splittings and coherence times.

In practice, the qubit frequency and the dephasing time, which are both tuned exponentially via the wire length and the coherence length, need to fall into an appropriate window for the effects discussed in this paper to be observable. If the wires are too long, the lifetime of the qubit may become limited by error processes not included here, for example, quasiparticle poisoning, which render the effects we discuss unobservable. Estimates for the quasiparticle poisoning times vary widely, but can easily exceed one microsecond [72]. Conversely, if the wires are too short, both the qubit precession period and its lifetime may be shorter than time-domain experiments can resolve. This timescale is mostly limited by how quickly the coupling between the qubit and the measurement dot can be tuned, which likely limits the resolvable timescales to about one nanosecond. Given these constraints, the ideal regime where the effects discussed in this paper can be observed still encompasses several orders of magnitude in timescale, and thus a significant window of wire lengths. As future direction, it would be interesting to develop protocols that can probe an even broader range of timescales.

Finally, we focused entirely on the minimal “tetron” qubit design for simplicity. A natural future direction is to extend our study to hexon devices that allow much more flexibility, including additional readout channels, measurement-based braiding, etc. Quantifying the role of noise in such higher-level applications would be an important and illuminating exercise. All in all, it is a remarkable feature of hardware-based topological qubits that we can directly simulate—in the time domain and with a single simulation—the effects of device-level noise on topologically protected quantum information.

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¹⁰Although for the qubit and noise model considered here, it is difficult to find a regime where clear “Ramsey oscillations” of $\langle i\gamma_1\gamma_3 \rangle$ persist and the basis choice gives rise to a clear difference.

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APPENDIX A: DETAILS OF NOISY TIME-DEPENDENT SIMULATIONS AND NOISE GENERATION

We perform time evolution following the Gaussian covariance matrix formalism described in Sec. III B. (See also Refs. [69,73]; for a self-contained overview, see Ref. [70].) Our numerical implementation is based on the Differential Equations.jl package in JULIA [74,75].

To generate a noise realization of some fluctuating quantity $\lambda(t)$ with a given noise power spectrum or correlation function, we use techniques based on fast Fourier transformation. This approach scales superlinear in the desired total time of the simulation (T_{sim} below). While for many types of noise (e.g., $1/\omega$ and $1/\omega^2$) more efficient ways of generating noise trajectories are known, in our context we never found this step to be a computational bottleneck. In addition, this approach has the advantage of being very general.

Consider a random variable $\lambda(t)$ with $\overline{\lambda(t)} = 0$ and $\overline{\lambda^2(t)} = (\delta\lambda_{\text{typ}})^2$, and correlation function $\overline{\lambda(t)\lambda(t')} = S(|t - t'|)$. At each time, the variable is drawn from a Gaussian distribution $P[\lambda(t)] \sim e^{-\lambda^2/2(\delta\lambda_{\text{typ}})^2}$. We can define the noise in frequency space $\lambda(\omega)$ as well as the noise power spectrum $S(\omega) \propto |\lambda(\omega)|^2$ as

$$\lambda(\omega) = \int_{-\infty}^{\infty} dt e^{-i\omega t} \lambda(t), \quad (\text{A1})$$

$$S(\omega) = \int_{-\infty}^{\infty} dt e^{-i\omega t} S(t). \quad (\text{A2})$$

For a simulation up to some finite time T_{sim} , we generate a discrete noise trajectory up to some time $T_{\text{noise}} \gg T_{\text{sim}}$ with a time step $\Delta t = T_{\text{noise}}/N$, i.e., the trajectory is defined on time slices $t_n = n\Delta t$, $n = 0, \dots, N-1$. It is important that $\Delta t \ll \omega_{\text{max}}$, where ω_{max} is the highest relevant frequency occurring in the noise. In the example of Gaussian noise characterized by a high-frequency cutoff κ , it is natural to set ω_{max} to some sufficiently large multiple of κ .

To numerically generate a single noise trajectory, we first create a white noise trajectory $Y(t_n)$ by drawing a sample on each time point from an independent and identical Gaussian ensemble of variance 1. We then perform a discrete Fourier transformation

$$Y(\omega_k) = \sum_{n=0}^{N-1} Y(t_n) e^{-i\omega_k t_n}, \quad \omega_k = k \frac{2\pi}{T_{\text{noise}}}. \quad (\text{A3})$$

We can then obtain our desired noise trajectory in Fourier space by taking

$$\lambda(\omega_k) = \left[\frac{S(\omega_k)}{\Delta t} \right]^{1/2} Y(\omega_k) \quad (\text{A4})$$

and finally performing an inverse Fourier transformation to obtain $\lambda(t_n)$:

$$\lambda(t_n) = \frac{1}{N} \sum_{k=0}^{N-1} \lambda(\omega_k) e^{i\omega_k t_n}. \quad (\text{A5})$$

In all these steps, it is convenient to choose N even so that the Fourier transformation is real. Furthermore, it is in many cases important to be able to sample $\lambda(t)$ for arbitrary t , for example, when integrating the Schrödinger equation with an adaptive time step. To this end, we perform a linear interpolation between time steps. If Δt is chosen sufficiently small, this will not incur significant numerical error.

In Fig. 12, we show the time dependence of all quantum-averaged quantities considered in this work for a single noise realization of our Ramsey-type protocol over one qubit precession period for the Kitaev tetron system of Secs. VI A and VII. We also show data necessary to derive the instantaneous qubit splitting [1]: the two lowest-lying instantaneous energies $\varepsilon_{1,2}$ and the determinant of the instantaneous orthogonal transformation U [see Eqs. (8) and (9)]. The left (right) panel corresponds to the initialization procedure based on the time-averaged (initial) Hamiltonian A_0 [$A(t=0)$] used in Sec. VI A (Sec. VII). Quantities with the subscript “fixed” denote measurement in a fixed basis throughout the evolution [e.g., $\gamma_i = \gamma_i^{(0)}$ in the left panel and $\gamma_i = \gamma_i(t=0)$ in the right panel], while quantities with the subscript “inst” denote measurement in the instantaneous basis derived from $A(t)$ [e.g., $\gamma_i = \gamma_i(t)$]. The data shown for $\langle i\gamma_1\gamma_3 \rangle_{\text{inst}} = \langle i\gamma_1(t)\gamma_3(t) \rangle$ are polluted by the numerical sign ambiguity described in Sec. IV B 3; we show it here over a single realization to illustrate this point. Comparing the envelope of $\langle i\gamma_1\gamma_3 \rangle_{\text{inst}}$ to $\langle i\gamma_1\gamma_3 \rangle_{\text{fixed}}$ demonstrates that for these parameters the choice of instantaneous versus fixed measurement basis does not have any impact on the results down to the level of each individual realization. Furthermore, comparing the dephasing diagnostics data in the top panels of Fig. 12 reveals, again for these chosen parameters, that the choice of initialization procedure makes negligible difference. For the leakage diagnostics in the bottom panel, the measurements are very close to unity, and we thus show an appropriately zoomed-in view. Note the observable high-frequency components in these measurements due to the small “quench” at $t=0$ when using the A_0 -based initialization procedure.

To calculate noise-averaged quantities, we run $N_{\text{real}} = O(10^3)$ independent such individual realizations. While this approach is relatively numerically demanding (compared, for example, to the master-equation-based approach of Ref. [61]), it has multiple advantages: For one, our techniques are completely general in terms of what noise models can be simulated and what physical quantities can be measured/noise averaged (in contrast to the methods of Ref. [61] which are more limiting in these respects); furthermore, our simulation strategy very closely mimics the actual experimental procedure we

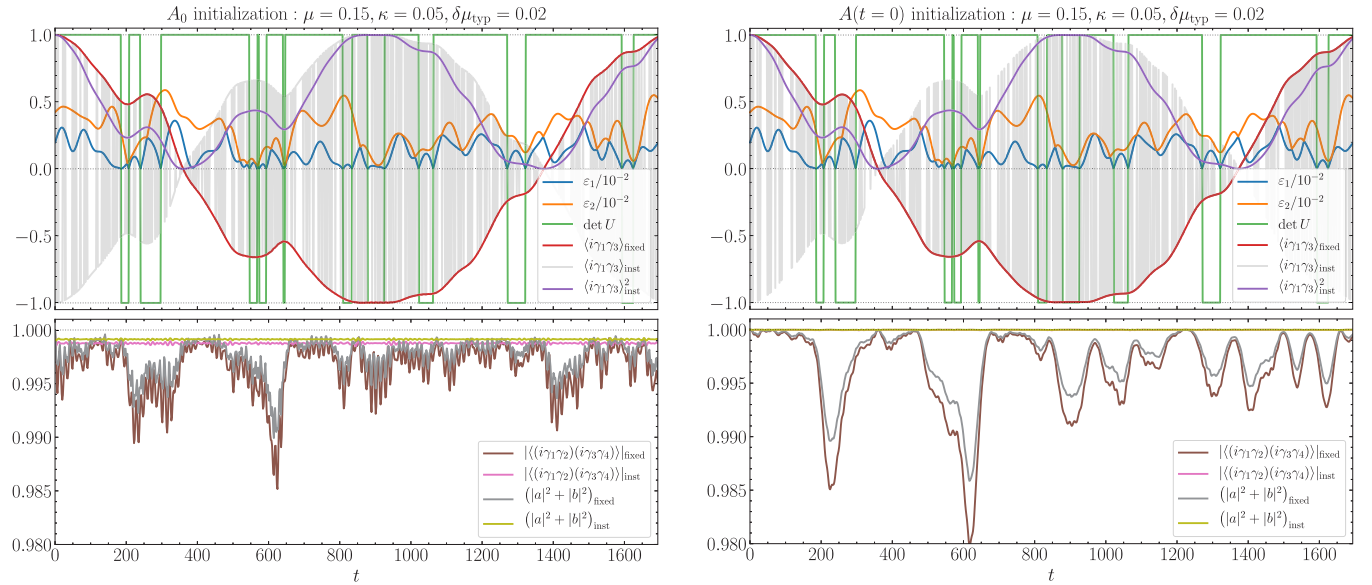


FIG. 12. Example time dependence of quantum-averaged quantities for a single noise realization our Ramsey-type protocol in the Kitaev tetron. Here, $\mu = 0.15$, $\kappa = 0.05$, and $\delta\mu_{\text{typ}} = 0.02$ with all other parameters and details identical to those in Secs. VIA and VII (see Figs. 3, 4, 5, and 10). In the left (right) panel, we show data obtained with the initialization procedure based on A_0 [$A(t=0)$] as in Sec. VIA (Sec. VII). The top panels include the dephasing diagnostics $\langle i\gamma_1\gamma_3 \rangle$ and $\langle i\gamma_1\gamma_3 \rangle^2$, as well as the (instantaneous) quantities $\varepsilon_{1,2}$ and $\det U$ from Eqs. (8) and (9); the bottom panels show the leakage diagnostics $\langle (i\gamma_1\gamma_2)(i\gamma_3\gamma_4) \rangle$ and $|a|^2 + |b|^2$. Subscripts in the legends indicate measurement in either the corresponding fixed or instantaneous (“inst”) basis.

propose (although in experiment noise and quantum averaging cannot be distinguished).

APPENDIX B: NOISE-AVERAGING ANALYSIS

Here, we provide details on the noise averaging of Eq. (33) assuming the Gaussian noise correlations specified in Eqs. (28)–(31). We will specifically evaluate $\exp(i \int_0^t dt' E(t'))$ with $E(t')$ given by the harmonic approximation in Eq. (26) and $\hbar = 1$ for notational simplicity; the noise average of $Q(t)$ follows straightforwardly from this quantity. It is convenient to discretize time (for intermediate stages of the calculation) and write

$$\overline{e^{i \int_0^t dt' E(t')}} = \frac{1}{Z} \int \mathcal{D}\lambda_i(t) e^{i\Delta t \sum_{i'=0}^t E(t')} \times e^{-\frac{1}{2} \sum_{i',t''} \sum_i S_i^{-1}(t'-t'') \lambda_i(t') \lambda_i(t'')}. \quad (\text{B1})$$

Here, Δt is the time interval used for discretization, the second line is the weighting factor that gives the desired noise correlations, and Z is a normalization defined as

$$Z = \int \mathcal{D}\lambda_i(t) e^{-\frac{1}{2} \sum_{i',t''} \sum_i S_i^{-1}(t'-t'') \lambda_i(t') \lambda_i(t'')}. \quad (\text{B2})$$

Next, we introduce

$$M_{ij}(t', t'') = M_{ij}^0(t', t'') + \delta M_{ij}(t', t''), \quad (\text{B3})$$

where

$$M_{ij}^0(t', t'') = \delta_{ij} S_i^{-1}(t' - t''), \quad (\text{B4})$$

$$\delta M_{ij}(t', t'') = \begin{cases} -i\Delta t \delta_{i',t''} E_{ij}''', & 0 < t' < t \\ 0, & \text{otherwise} \end{cases} \quad (\text{B5})$$

and also

$$v_i(t') = \begin{cases} i\Delta t E_i', & 0 < t' < t \\ 0, & \text{otherwise.} \end{cases} \quad (\text{B6})$$

These definitions allow us to write

$$\overline{e^{i \int_0^t dt' E(t')}} = \frac{e^{iE_0 t}}{Z} \int \mathcal{D}\lambda_i(t) e^{\sum_i \sum_{t'} v_i(t') \lambda_i(t')} \times e^{-\frac{1}{2} \sum_{ij} \sum_{t',t''} \lambda_i(t') M_{ij}(t', t'') \lambda_j(t'')}. \quad (\text{B7})$$

Note that the t', t'' sums are unrestricted above. We can now perform the Gaussian integration to obtain

$$\overline{e^{i \int_0^t dt' E(t')}} = e^{iE_0 t} e^{-\frac{1}{2} \ln(\frac{\det M}{\det M^0})} \times e^{\frac{1}{2} \sum_{ij} \sum_{t',t''} v_i(t') M_{ij}^{-1}(t', t'') v_j(t'')}. \quad (\text{B8})$$

To proceed we first expand the log term in Eq. (B8) to second order in δM :

$$\begin{aligned} \ln \left(\frac{\det M}{\det M^0} \right) &= \ln \det [I + (M^0)^{-1} \delta M] \\ &\approx \text{Tr}[(M^0)^{-1} \delta M] - \frac{1}{2} \text{Tr}[(M^0)^{-1} \delta M]^2 \\ &= -it \sum_i D_i^2 E_{ii}'' \\ &\quad + \frac{1}{2} \sum_{ij} (E_{ij}'')^2 \int_0^t dt' dt'' S_i(t' - t'') S_j(t'' - t'). \end{aligned} \quad (\text{B9})$$

The third line above encodes the leading shift in the qubit precession frequency [cf. Eq. (32)]. At long times, $t \gg \tau_i$, the

integrals in the fourth line become

$$\int_0^t dt' dt'' S_i(t' - t'') S_j(t'' - t') \approx t \frac{2\sqrt{\pi} \tau_i \tau_j}{\sqrt{\tau_i^2 + \tau_j^2}} (D_i D_j)^2. \quad (\text{B10})$$

In the second line of Eq. (B8) we simply replace $M^{-1} \approx (M^0)^{-1}$; the next-leading correction provides a higher-order shift to the qubit precession frequency compared to that captured above. This approximation yields

$$\begin{aligned} \sum_{ij} \sum_{t't''} v_i(t') M_{ij}^{-1}(t', t'') v_j(t'') \\ \approx - \sum_i (E'_i)^2 \int_0^t dt' dt'' S_i(t' - t'') \approx t 2\sqrt{\pi} \sum_i \tau_i (D_i E'_i)^2, \end{aligned} \quad (\text{B11})$$

where on the far right side we again assumed the long-time limit.

Putting everything together, we find

$$\begin{aligned} \overline{e^{i \int_0^t dt' E(t')}} = e^{i(E_0 + \frac{1}{2} \sum_i D_i^2 E''_{ii})t} \\ \times e^{-\sqrt{\pi} \left[\sum_i \tau_i (D_i E'_i)^2 + \frac{1}{2} \sum_{ij} \frac{\tau_i \tau_j}{\sqrt{\tau_i^2 + \tau_j^2}} (D_i E''_{ij} D_j)^2 \right] t} \end{aligned} \quad (\text{B12})$$

at $t \gg \tau_i$. Upon restoring explicit \hbar 's, this result indeed recovers the qubit precession frequency and dephasing time quoted in Eqs. (32) and (35).

The short-time limit $t \ll \tau_i$ can be easily treated as well. Here, we can simply write

$$\int_0^t dt' dt'' S_i(t' - t'') S_j(t'' - t') \approx t^2 S_i(0) S_j(0) = (D_i D_j t)^2, \quad (\text{B13})$$

$$\int_0^t dt' dt'' S_i(t' - t'') \approx t^2 S_i(0) = (D_i t)^2 \quad (\text{B14})$$

in Eqs. (B10) and (B11), so that

$$\begin{aligned} \overline{e^{i \int_0^t dt' E(t')}} = e^{i(E_0 + \frac{1}{2} \sum_i D_i^2 E''_{ii})t} \\ \times e^{-\frac{1}{2} [\sum_i (D_i E'_i)^2 + \frac{1}{2} \sum_{ij} (D_i E''_{ij} D_j)^2] t^2}. \end{aligned} \quad (\text{B15})$$

The noise-averaged $Q(t)$ is then

$$\overline{Q(t \ll \tau_i)} \approx \cos(\omega_0 t) e^{-(t/T_2^{\text{short time}})^2} \quad (\text{B16})$$

with (restoring \hbar 's)

$$T_2^{\text{short time}} = 2\hbar^2 \left[\sum_i (D_i E'_i)^2 + \frac{1}{2} \sum_{ij} (D_i E''_{ij} D_j)^2 \right]^{-1}. \quad (\text{B17})$$

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